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(54) Title: MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

(57) Abstract

Assay methods for determining whether a peptide is likely to be immunogenic are based on a computer modeling of binding to a Class II MHC DR1 receptor. This is confirmed by competitive inhibition binding assays. The peptides are useful for eliciting an immune response for vaccination or the production of antibodies or T-cells.

Applicants: Moses Rodriguez and Daren Ure

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Exhibit 12

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MODEL FOR TESTING IMMUNOGENICITY OF PEPTIDES

Government Interest

The invention described herein may be manufactured, licensed and used by or for governmental purposes without the payment of any royalties to us thereon.

Cross Reference

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This application is a continuation-in-part of U.S. Patent application Serial No. 08/064,559, filed May 21, 1993, and the present application incorporates U.S. Patent Application Serial No. 08/064,559 in its entirety by reference.

15 Field of the Invention:

This invention relates to a means of predicting potential of a peptide for eliciting immune response.

Background of the Invention:

Among the numerous steps required for an immunological response to occur is the presentation of the antigen by macrophages to the B-cell or T-cell. This presentation is mediated by the Class I and Class II major histocompatibility complex (MHC) molecules on the surface of the cell. The MHC molecules hold antigens in the form of the peptide fragments and together with the receptor molecule on the T-cells, form a macromolecular complex that induces a response in the T-cell. Therefore, a necessary step in an immune response is the binding of the antigen to the MHC.

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Recent single crystal X-ray structures of human and murine Class I MHC's have been reported. Analysis of these crystal structures have shown that antigenic peptides lie in the so-called binding cleft for presentation to the T-cell. This cleft is formed by α_1 and α_2 domains and by eta-strands from each domain forming the floor. Furthermore, the sequence polymorphism among Class I molecules can result in alterations of the surface of the cleft forming different pockets. Peptide side chains may insert into these pockets. Thus, different pockets may interact with different side chains. This implies the mechanism for the peptide specificity of Class I MHC's. Peptides bound to the Class I MHC's in the crystal structures were found to have both the amino and carboxy termini tightly held by the MHC. There were few interactions near the middle of the cleft. Hence the bound peptide is allowed to bend slightly in the center. observed binding mode helped to explain the apparent partial specificity of peptide sequence and the allowed variation in peptide length found among peptides isolated from Class I MHC's.

The precise mode of binding of peptides to Class II MHC molecules is less clear. While a single crystal X-ray diffraction structure for the HLA-DR1 MHC has been shown, the coordinates have remained unavailable. However, currently available theoretical and experimental results help form a hypothesis that the binding of a peptide to Class II MHC is similar to that observed with Class I. First, it is noted that the Class II binding cleft is structurally similar to

that of Class I. This was concluded based upon a sequence analysis of 26 Class I and 54 Class II amino acid sequences.

Unlike with Class I molecules, self-peptides isolated from murine I-A^b and I-E^b, from murine I-A^d and from human HLA-DR1 molecules were found to be varied in size (13 to 25 residues long). The peptides isolated from the murine I-A^b and I-E^b molecules had heterogenous carboxy termini while those from I-A^d and HLA-DR1 had ragged termini at both ends. The varying lengths indicate that the amino and carboxy termini of the peptides were not critical for the binding. One or both termini may protrude from the binding site and be available for further processing. The residues critical for binding were proposed to be at the ends of the peptide as opposed to the center.

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Summary of the Invention:

It is the purpose of this invention to provide a method for preliminary screening of peptides for ability to elicit an immune response. Structural homology techniques were used to model a receptor (the Class II MHC is exemplified). This model makes it possible to preliminarily screen peptides for antigenic properties. By modifying the peptide to "fit" into the receptor it is possible to identify methods of rendering non-immunogenic peptides immunogenic.

The preliminary screening of peptides for immunogenicity comprises the steps of (1) creating a molecular model of a receptor followed by minimizing the model created, 2) modeling a peptide to be tested and minimizing the model of the peptide, then testing the

fit of the model of the peptide into the model of the receptor to produce a composite minimized receptor/minimized peptide model.

Upon finding an acceptable fit, the peptide may then be screened by a binding assay for actual binding to Class II MHC as a further test for immunogenicity.

It has been found that when the model of the peptide can not i fitted into the model of the receptor, the peptide will lack immunogenicity. While not all peptide models which can be made to "fit" into to model of the receptor will be effective as immunogens the screening methods of the invention may make it possible to avoi undue biological testing of inappropriate peptides. By using the model, it is also possible to alter peptides to accommodate the receptor. Hence, the invention has both predictive and drug design applications.

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Brief Description of the Figures:

Fig. 1 shows the HLA-aw68 $lpha_1$ and $lpha_2$ domains with DR1 $lpha_1$ and eta_1 domains.

Figs. 2-30 are a printout of the minimized coordinates of the 20 receptor.

Figs. 31 and 32 shows the effects of various peptides inhibiting the binding of labeled hemagglutinin in a competitive binding assay.

25 <u>Detailed Description of the Invention:</u>

In order to understand and better predict peptide interaction with Class II MHC's and as an aid for synthetic peptide vaccine design, a structural homology model of HLA-DR1 molecule was made

using the Class I HLA-aw68 as a reference molecule. For purposes c: this analysis, numerous conserved residues were aligned leading to a proposed three-dimensional model for the Class II structure very similar to that of Class I. This model retained the overall conformation of a Class I MHC and agreed with a considerable amount of the published data. Furthermore, peptides shown to bind to DRI were docked in the binding cleft of the model and analyzed. The results agree with the experimental binding data presented here. Hence, it is shown that the structural homology model reported here is useful for screening Class II MHC functionality.

It had been hypothesized that few peptide residues may be required for binding to DR1. By substituting residues into the influenza hemagglutinin 307-319 T-cell epitope (HA) it had been determined that a single tyrosine at 308 was required for binding. A synthetic peptide with the tyrosine at position 308 and a lysine at 315 was found to bind DR1 as well as the native peptide. Hence, it was concluded that few peptide residues determine the high affinity binding to DR1.

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The peptides produced according to the present invention may be used alone or chemically bound to another peptide and/or carrier in order to elicit an immune response. An immune response is elicited by administering a peptide to an animal in an effective dose and by an effective route of administration. Typically the peptide will be administered with an immunologically acceptable carrier. The route of administration, dosages, times between multiple administrations will be based on the particular peptide and are standard operations of those skilled in the art.

of particular interest are peptides from pathogenic microorganisms and neoplasms. In such an example, a vaccine may be formed with the peptide and any known immunological carrier and may be administered prophylactically or therapeutically. The immune response may be elicited for a number of reasons other than for prophylaxis or therapy such as increasing antibody production for the harvesting of antibodies, or increasing specific B-cell or T-cell concentration for the production of hybridomas or cellular therapy.

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The choice of host animals is limited only to those capable of an immune response. Preferred hosts are mammals, more preferred are humans.

The vaccine may contain plural peptides with each peptide corresponding to the same or different antigens. The peptides may be used unbound or they may be chemically bound to another peptide or an unrelated protein or other molecule. A preferred vaccine preparation contains a plurality of peptides chemically bound to a larger more immunogenic peptide.

The peptide may be adsorbed, bound or encapsulated in a biodegradeable microsphere, microcapsule, larger carrier or a combination of these. The carrier may have a slow or controlled release property thereby releasing the peptide under appropriate conditions and times for enhanced immunization. This is particularly important when administering the peptide orally where stomach acid can degrade the peptide.

Another embodiment of the present invention is to modify the amino acid sequence of a peptide to enhance its immunogenicity.

This is done by modifying the natural peptide sequence to bind to

the Class II MHC receptor DR1 with superior binding affinity for a Class II MHC receptor DR1 than the natural peptide sequence. This modified peptide is considered a synthetic peptide. Alternatively, the sequence may be modified to have a greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1.

Many amino acid changes are acceptable in the formation of a synthetic peptide. The changes may be for similar types of amino acids such as leucine for isoleucine or they may be for diverse types such as tyrosine for lysine.

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Materials and Methods:

The structural homology model for the DR1 Class II MHC was constructed using the QUANTA molecular modeling package (vision 3.2, Molecular Simulations, Inc., Burlington, MA) with the CHARMM and Protein Design modules. After alignment of the sequences as described below, gaps and loops were energy minimized using 100 steps of steepest descents minimization followed by 100 steps of adopted basis set Newton-Rapheson (ABNR) minimization. Large gaps were closed using a fragment database from a selected set of high-resolution crystal structures. The resulting structure was minimized in vacuo using 1000 steps of steepest descents followed b an additional 1000 steps of ABNR minimization. A distance related electrostatic function was used in all calculations with a dielectric constant of 1.0. Non-bound parameter lists were updated every 20 steps with a cutoff distance of 15.0Å. Non-bonded calculations were performed using a shifted potential function between 11.0Å and 14.0Å. An extended atom set was used with only

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polar hydrogen atoms specifically placed. There were no explicit hydrogen bond energy calculations performed.

All peptides were initially modeled using QUANTA in an extende chain conformation and subjected to 500 steps of ABNR minimization. The resulting structures remained essentially in extended chain conformations. Individual peptides were manually docked in several different orientations into the binding cleft region of the minimized DR1 structure. The resulting bimolecular complex was subjected to 5000 steps of steepest descents minimization with non-bonded interactions updated every five steps. After minimization, bound peptides remained essentially in extended chain conformations. The lowest energy complexes for each peptide were selected for further analysis.

The selected peptide and DR1 complexes and the minimized DR1 model were subjected to the following molecular dynamics regimen: 300 steps of heating to 300°K, 600 steps of equilibration at 300°K, and 1100 steps of production dynamics. During this simulation, the DR1 Ca atoms were constrained in their starting positions. All non-bonded interaction parameters were as stated for the minimization procedure. The lowest energy structure during the course of the production dynamics was selected and subjected to the 5000 step minimization procedure described previously with the Ca restraints removed. The resulting structures were used for the binding energy calculations and for hydrogen bonding analysis.

Hydrogen bonds were determined using the QUANTA default parameters. Maximum allowed distances were 2.5Å between a hydrogen and the acceptor atom and 3.3Å between the donor and acceptor atom

The minimum angle allowed between any set of atoms forming a hydrogen bond was 90°.

Competitive Inhibition Binding Assay:

HA peptide (the influenza hemagglutinin 307-319 T-cell epitope) was labeled with ¹²⁵I. The labeled HA peptides were then allowed to interact with purified DR1 molecules during incubation to allow formation of peptide/DR1 complexes. After incubation, the peptide/DR1 composition was exposed to a native gel for chromatographic separation or passed through a spun column to separate labeled peptide/DR1 complex and free labelled peptide. When unlabeled peptides were added before incubation of labeled HA peptides and DR1, and if the unlabelled peptides had capacity for binding to DR1 simultaneous with ¹²⁵I-HA, there was a resultant decrease in radioactive signal associated with the DR1. The extent of this decrease directly related to the binding capacity of the unlabeled unknown peptide.

Structural Homology Model for the DR1 Molecule:

The structural homology model was created, the reference molecule being the crystal structure of HLA-aw68. The HLA-aw68 coordinates and subsequent sequence were obtained from the entry 2HLA in the Brookhaven Protein Data Bank released January 15, 1991, which is incorporated herein by reference. The sequence for the DR1 molecule was for the α_1 domain was reported by Klein and for the β_1 domain, the study reported by Todd et al. (Nature 329, 599 (1987)).

The sequence alignment is based on Brown et al. (Nature 332, 845 (1988)). The complete alignment and numbering scheme for both

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are seen in Figure 1. The Class II, β , and Class I α , domains regions were conserved with some variations at the ends where the two MHC's have different loop regions. The fourth B-strand in the lphadomain of HLA-aw68 (residues 30-38) is disrupted in the DR1 model. Only three residues are in a β -sheet conformation, probably due to the inserted glycine at position 28 before the strand and the large deletion in the loop region immediately after the strand. The two alpha-helical regions are clearly maintained. Both helices have been observed to be discontinuous in the Class I molecules and are similar in the DR1 model. The α_i domain helix is long and curves from residues 49α to 76α without significant disruption. essentially a single continuous helix. However, the $lpha_2$ helical region is broken into two separate helices as with the Class I molecules. A short helix (52-63) is separated from a longer helix (68-94) by a deformed region without secondary structure. deformation is more pronounced in the DR1 model as opposed to the Class I molecules due to an insertion.

Influenza Hemagglutinin Peptide with DR1:

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The amino acid residues 307-319 of influenza hemagglutinin (Pro-Lys- Tyr-Val-Lys-Gln-Asn-Thr-Leu-Lys-Leu-Ala-Thr) make up a well-documented linear T-cell epitope which has been shown to be HLA-DR1 restricted. With the demonstration that the influenza hemagglutinin epitope (referred to as the HA peptide) binds DR1,it was chosen to be modeled into the binding cleft.

The peptide was initially inserted into the cleft so that Leu 11 HA was in the vicinity of the hydrophobic pocket. This allowed Asn 7 to be near the middle charged and polar groups of the cleft.

The remaining residue of the motif (Lys 2) was near the vicinity of the remaining charged and polar residues at the end of the cleft.

The only adjustment to the starting conformation was a slight rearrangement of the terminal peptide proline and Tyr 3 to alleviate obvious bad contacts.

After the energy minimization of the bimolecular complex, the total energy was reduced to 483 kcal/mol. This reduction in energy was accomplished by alleviation of several bad contacts and also be formation of several hydrogen bonds. The sticking feature of this mode is lack of hydrogen bonds in the carboxy terminal half of the peptide. Only one hydrogen bond is identified between the backbone carbonyl group of Leu 9 and the side chain of the β_1 Asn 77. In contrast, the amino terminal half has eleven identified interactions. Four of these interaction involve the peptide backbone residues Tyr 3, Val 4, and Gln 6. The remainder involve the side chains of Lys 2, Tyr 3, Lys 5 and Gln 6. Interestingly, Lys 5 is involved in more interactions (three) than Lys 2 (only 2). No interactions were observed as anticipated with Asn 7. Instead, it was the glutamine at position 6 donating a hydrogen bond to the ϵ Asn 62. No interactions were observed for the amino and carboxy termini.

HA-YK Peptide with DR1:

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The binding of the HA-YK peptide (Ala-Ala-Tyr-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Ala-Lys-Ala-Ala) to the DR1 model was tested. In aligning the peptide in the cleft, it was deemed logical to insert the tyrosine residue into the hydrophobic region of the binding cleft. The lysine would then be in position to interact with the

hydrophilic groups in the other half of the cleft. The resulting peptide orientation is the opposite that used for the HA and the CS3 (defined below) peptides. With the peptide oriented as described, the final docking position for the peptide was unclear. The hydrophobic pocket is quite large, and, at least in this model, could accommodate the peptide tyrosine in a number of positions by sliding the peptide lengthwise through the cleft. However, repositioning the peptide also repositions the lysine. There were primarily two positions for the lysine: one with the lysine inside the cleft and the second with it outside. Of the two positions, the former was the lower in energy by 46 kcal/mol and had the greater number of interactions with the protein (11 vs. 7). Thus, the preferred orientation of the peptide appears to be with the lysine inside the binding cleft region.

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CS3 subunit Pilin Peptide with DR1:

The suspected T-cell epitope for CS3 pilus subunit 63-78

(Ser-Lys-Asn-Gly-Thr-Val-Thr-Trp-Ala-His-Glu-Thr-Asn-Asn-Ser-Ala)

was modeled with the DR1 molecule. The peptide was inserted with

lysine inside the cleft in the hydrophilic region. This placed the

Thr 5 in the center of the binding cleft and the tryptophane

(residue 8) near the hydrophobic region. The resulting minimized

model had ten interactions between the peptide and the protein,

three interactions with the peptide backbone and five with the

peptide side chains. The remaining two were with the amino termina

of the peptide. All of the interactions were in either the first

three residues, His 10 or Glu 11 in the peptide. No interactions

were observed in the center of the cleft or residues four through nine.

CFA/1 with DR1:

A peptide identified as CFA/1 (colonization factor antigen)

(Val-Gly-Lys-Asn-Ile-Thr-Val-Thr-Ala-Ser-Val-Asp-Pro) was prepared

and an attempt was made to "fit" the molecule into the cleft of the

DR1. The lysine at position 3 prevented insertion of the peptide.

10 Results:

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The peptides chosen to dock in the DR1 model are shown in Table

1. The peptides were docked manually in several orientations into
the DR1 model. The peptides were then tested in biological binding
assays with the following results:

Table I

Peptide	Molecular Model predicted binding	Binding in the bioassay
HA (influenza hemagglutinin)	Yes	Yes
HA-YK (synthetic peptide)	Yes	Yes
CS3 Pilin subunit	Yes	Yes
CFA/1	No	No

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Quantitative measurement of the inhibition of CS3 63-78 and H^2 306-318 as compared to controls is shown in Fig. 31.

The binding energy was calculated as the difference between the final DR1 and peptide complex and the sum of the energies for the minimized DR and peptide models individually. The data is shown in Table II.

Table II.

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Peptide	Protein	Residues	Sequence	Binding Energy (kcal/mol)
НА	Influenza hemagglutinin	306-318	PKYVKQNTLKLAT	-283
HA-YK	synthetic peptide		ААҮАААААКАА	-216
CS3	CS3 pilin subunit	63-78	SKNGTVTWAHETNNSA	-245

CS6α and CS6ß with DR1

Colonization factor antigen IV (CFA/IV is an antigen on the

surface of many enterotoxigenic *E. coli* one component of which is

CS6. CS6 has two major subunits and a number of minor subunits.

Several peptides from CS6 have been sequenced and assayed for
potential inhibition of radiolabeled HA (306-318)/DR1 complex as a
measure of immunogenicity. The sequences of the subunits are shown
in Table III.

Table III.

Peptide_	Amino Acid Residues	Sequence
CS6α6	63-75	DEYGLGRLVNTAD
CS6α7	80-92	IIYQIVDEKGKKK
CS6α8	111-123	LNYTSGEKKISPG
CS6ß1	3-15	WQYKSLDVNVNIE
CS6ß2	42-54	QLYTVEMTIPAGV
CS6ß3	112-124	TSYTFSAIYTGGE
CS6ß4	123-135	GEYPNSGYSSGTY
CS685	133-145	GTYAGHLTVSFYS

These peptides were assayed for inhibition of radioactively labeled HA(306-318)/DR1. The results are demonstrated in Fig. 32.

The foregoing description of the specific embodiments reveal the general nature of the invention so that others can, by applying current knowledge, readily modify and/or adapt for various applications such specific embodiments without departing from the generic concept, and, therefore, such adaptations and modifications should and are intended to be comprehended within the meaning and range of equivalents of the disclosed embodiments. It is to be understood that the phraseology or terminology employed herein is for the purpose of description and not of limitation.

All references mentioned in this application are incorporated by reference.

We Claim:

- 1. A method of preliminarily screening peptides for immunogenicity comprising the steps of:
- 1) creating a molecular model of receptor DR1 Class II MHC and minimizing the model of the DR1;
 - 2) modeling a peptide to be tested and minimizing the model of the peptide; and
- 3) testing fit of model obtained in step 2 into the model
 10 obtained in step 1 to produce a composite receptor/peptide model.
 - 2. A computerized model comprising a model of the DR1 molecule having fitted in a cleft therein a model of a peptide.
- 3. A method of claim 1 wherein, additionally, the receptor/peptid model is subjected to computer-simulated heating.
 - 4. A method of claim 1 further comprising, assaying the peptide becompetitive inhibition binding to a Class II MHC receptor DR1.
 - 5. A minimized peptide capable of binding to a Class II MHC receptor DR1 and inhibiting the binding of HA (306-318).
- 6. A synthetic peptide, wherein the amino acid sequence of the
 minimized peptide according to claim 5 has been modified to have a
 superior binding affinity for a Class II MHC receptor DR1 to form;
 least a portion of the synthetic peptide.

7. A synthetic peptide, wherein the amino acid sequence of the minimized peptide according to claim 5, has been modified to have greater inhibition of HA (306-318) binding to a Class II MHC receptor DR1 to form at least a portion of the synthetic peptide.

- 8. A synthetic peptide according to claim 6, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 9. A synthetic peptide according to claim 7, wherein an amino acid has been modified from a charged amino acid to an uncharged amino acid.
- 10. A synthetic peptide according to claim 8, wherein saiduncharged amino acid is alanine.
 - 11. A synthetic peptide according to claim 9, wherein said uncharged amino acid is alanine.
- 12. A minimized peptide according to claim 5, wherein the sequence is selected from the group consisting of PKYVKONTLKLAT, AAYAAAAAKA; and SKNGTVTWAHETNNSA.
- 13. A minimized peptide according to claim 5, wherein the sequence is contained in a CFA.

14. A minimized peptide according to claim 13, wherein the sequence is selected from the group consisting of DEYGLGRLVNTAD, IIYQIVDEKGKKK, LNYTSGEKKISPG, WQYKSLDVNVNIE, QLYTVEMTIPAGV, TSYTFSAIYTGGE, GEYPNSGYSSGTY and GTYAGHLTVSFYS.

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- 15. A vaccine comprising:
 - a minimized peptide according to claim 5; and an immunologically acceptable carrier.
- 10 16. A vaccine comprising:
 - a synthetic peptide according to claim 6; and an immunologically acceptable carrier.
 - 17. A vaccine comprising:
- a synthetic peptide according to claim 7; and an immunologically acceptable carrier.
 - 18. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 15.
 - 19. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 16.

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20. A method of eliciting an immune response in an animal comprising administering said animal with the vaccine according to claim 17.

8904	-i -	CSUSPRITE 9		GRG 19 EPRPIAVGYV 29		SDAASQRHEP 48 1	rapuleqegp etvurleppg
DR1 .,	E -	TKEEÜVIIQA 11	EFYLII FUQ 19	SG EFFEOR 27			•
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DR1 a,	S0 R	PASFEAGGA 60	LAHTAVDKAH 70	LEIMTKÄSIIY 80	TPI Chrange 8	LWOLKPECHIF 18	FIICTERVRLL
OR1 B,				-			
8920	116	RQDAYDGKDY 12	114 RQDAYDGKDY 124 1ALKEDLRSW 134 TAADMAA QT 143 TKUXWEAA H 152 VAEGURAYLE 162 GTCVEWLRR"	(((((((()	CCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCC	((* bollx)))) 2 VAEQURAYLE 162	GTCVEWLRR:
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DR1 B,	28	ERCIYNQEES 31	0 VIJEDSDVGEY 40	RAVTELCRIO 58	ALI WISHING		
B9) r	172	LENGKETLOR					
DK1 .							
DR1 B,	87	esftvqrrvu					

FIG. 1

Conserved residues Polyworphic residues

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1639								0.00000
1	1 ILE		-53.41835	5 -52.87964	96.86949 96.22549		1 .	0.00000
2 3	1 ILE 1 ILE	HT1 UT2	-53 40505	3.37379 553.33354	96.22349		ī	0.00000
4	1 ILE	HT3	-53.81151	-52.85195	97.84341		1	0.00000
5	1 ILE	CA		-51.45945			1	0.00000
6	1 ILE	CB		5 -51.09296		λl	1	0.00000
7	1 ILE	CG2		7 -51.39510			1	0.00000
8	1 ILE	CG1		-49.65164			1	0.00000
9	1 ILE	CD		-49.3065			1	0.00000
10	1 ILE 1 ILE	0		5 -50.79352 2 -51.45486			1	0.00000
11 12	2 LYS	ห		-49.57271			2	0.00000
13	2 LYS	11		-49.07042	97.19065	Al .	2	0.00000
14	2 LYS	Cλ		-48.82990			2	0.00000
15	· 2 7.YS	CB	-51.48674	-49.22996	100.05168	N1	2 2	0.00000
16	2 LYS		51.65942	-50.46122 -50.76541		A1	2	0.00000
17 18	2 LYS 2 LYS	CD CE		-51.67024			2	0.00000
. 19	,2 LYS	NZ		51.62033			2	0.00000
20	2 LYS	1121		-52.15413	104.71032	λί	2	0.00000
21	2. LYS	HZ2	-48.66152	-52.03086	103.36182		2	0.00000
~22	2 LYS				104.08530		2	0.00000
23	2 LYS	С		-47.37619			2	0.00000
. 24	2 LYS	0		-47.08993 -46.48610			3	0.00000
25 26	3 GLU	H H	-54.93373	-46.40010	100.69754		3	0.00000
27	3 GTO	CA		-45.05669	99.49342	Al	3	0.00000
28	3 GTA	CB		-44.67728	98.75869	λl	3	0.00000
29	3 GTA	CG		-43.30721			3	0.00000
30	3 CIU	CD		-43.14982			3 3	0.00000
31	3 GLU	021		-42.20412 -43.96983	96.38367 97.23742		3	0.00000
32 33	3 GLU	OE2 C	-50.30000	-44 3058B	100.80334		3	0.00000
34	3 GLU	o		-44.80289		Al	3	0.00000
35	4 GLU	N			100.72198		4	0.00000
36	4 GLU	H	-51.89913	-42.70887	99.83885		4	0.00000
37	4 GLU	CA	-51.71490	-42.41569	101.93532	Al	4	0.00000
38	4 GLU	CB.	-50.23606	-42.65775	102.23912	ri Li	4	0.00000
39 40	4 GLU 4 GLU	CD	-49.88908	-44.20822	102.86978	Al	4	0.00000
41	4 GLU	0Z 1	-47.71593	-43.20739	103.12446	Al	4	0.00000
52	4 GLU	OE 2	-47.87485	-45.31826	102.72475	Al .	4	0.00000
43	4 GLU	С	-51.85859	-40.92476	101.75610	A1	4	0.00000
44	4 GLU	0	-51.85445	-40.40438	100.64776	Al Ni	4 5	0.00000
45 46	5 HIS 5 HIS	и н .	-51.98/58	-40.25490	102.89941	Al	5	0.00000
47	5 HIS	H. Ca	-52.02510	-38.79739	102.88794	スコ	5	0.00000
48	5 HIS	CB	-52.95268	-38.30654	104.00423	Al	5	0.00000
49	5 HIS	CG	-54.39292	-38.58657	103.64511	Al	5	0.00000
50	5 HIS	NDI	-55.01336	-38.04007	102.58831	λl	5 .	0.00000
51	5 HIS	HD1	-54.63216	-37.40030	101.93314	A1	5 5	0.00000
52 53	5 HIS	CD2	-55.29163 -56.46563	-39.42491 -30.33333	104.31043	Al	5	0.00000
53 54	5 KIS 5 KIS	NE2 CE1	-56.29489	-38.51954	102.57197	<u> </u>	5	0.00000
55	5 HIS	C	-50.64149	-36.20241	103.06558	λl	5	0.00000
5.6	5 HIS	0	-49.7590E	-36.76174	103.65940	Al	5	0.00000 0.00000
57	6 VAL	1: :-	-50.46014	-37.02655	102.46317	A1	6 6	0.00000
56	6 VAL	H C:	-51.22869 -49.12695	-36.576	102.46428	£1	6	0.00000
5 9 6 0	6 VAL 6 VAL	CA CB	-48.60121	-36.33669	101.01420	A 2	£	0.0000
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62	6 VA:			4 -37.4664			0.00000
€3	6 VAL	. с		5 -35.0019			0.00000
64	6 VAI				1 102.87457		0.00000
65	7 115		-48.1152	7 -34.5212	103.65431		0.00000
66 67	7 ILE 7 ILE		-47.3597	2 -35.13140 6 -33 0869	3 103.91409 7 103.98819		0.00000
68	7 ILE		-48.0930 -48.6919	7 -32.8681	3 105.39701	λ1 7	0.00000
69	7 ILE		-47.9632	2 -33.7131	7 106.43001		0.00000
70	7 ILE			6 -31.3958			0.00000
71	7 ILE	CD	-49.2884	6 -31.2048	107.23523		0.00000
72	7 ILE	С	-46.6938	1 -32.5011	103.87753		0.00000
73	7 ILE	0		5 -33.10109			0.00000
74	8 ILE	И		-31.32789			0.00000
75	8 ILE	H		3 -30.79777 5 -30.85113			0.00000
76 77	8 ILE	CA CB		-31.21426			0.00000
78	8 ILE	CG2		-30.98133			0.00000
79	8 ILE	CG1	-44.01583	-30.53590	100.51961	A1 8	0.00000
80	8 ILE	.:CD			101.12422		0.00000
81	8 ILE	С			103.03239		0.00000
82	8 ILE	· O			102.63903		0.00000
83	9 GLN	N			103.70229		0.00000
84	9 GLN	н			104.05676		0.00000
85	9 GLN	CA	-43.27910	-27.85712	103.71549 2	Al 9 Al 9	0.00000
£5 87	9 GLN	CB ·	-44.97213	-27.43231	105.94550	A1 9	0.00000
85	9 GLŅ	CD	-43.92932	-26.69043	107.36359	A1 9	0.00000
89	9 GLN	OE1	-44.59992	-27.05224	108.31811	A1 9	0.00000
90	9 GLN	NE2	-42.89278	-25.86874	107.50418	A1 9	0.00000
91	a Crn	HE21	-42.31986	-25.58069	106.73881		0.00000
92	a CIN	HE22			108.41526		0.00000 0.00000
93 94	9 GLN	C	-42.00840	-27.79728 -28.56703	102.89330 /	_	0.00000
95	9 GLN 10 ALA	0 N		-26.84230	·		
96	10 ALA	H		-26.24022	101.81084		
97	10 ALA	CA		-26.60034	101.15833 2		0.00000
98	10 ALA	СВ	-41.09424	-26.95993	99.69281 3	1 10	
99	10 ALA	С	-40.41733	-25.14834	101.25674 3		
100	10 ALA	0		-24.24680	101.50542		
101	11 Gro	Ŋ	-39,11637	-24.95043	101.08226		
102 203	II GLU	H CA		-23.61077	100.86462 7		
104		·CB			102.46404 8		
105	11 GLU	CG	-38.25740	-24.15021	103.73032 A	11	0.00000
106	lî GLU	CD	-37.31724	-24.35236	104.86668 A	11	
107	11 CLD	OEl	-37.72498	-24.91514	105.87939 A	1 11	
108	11 CTO	OE2			104.74301 8		
109	11 GLU	С			100.04756 A 99.37097 A	1 11	0.00000
111	11 GLU 12 PHE	0 N	-37.17390 -37.8688	-21.87280	.99.78525 A	1 12	
112	12 PKE	Н		-21.25141			
113	12 PHE	CA	-37.20863		98.59813 A		0.00000
114	12 PHE	CB	-38.26225	-21.06791	97.51950 A		0.00000
115	12 PHE	CG	-37.93682		96.27668 A		0.00000
116	12 PHE	CD1	-38.35291	-23.20899	96.17465 A		0.00000
117	12 PHE		-37.21678		95.22261 A		0.0000
118 119	12 PHE		-38.04712 -36.91098		95.01356 A 94.06122 A		0.00000
120	12 PHE		-30.91096		93.96022 A		0.00000
121	12 PHE		-36.45949		98.90209 %		0.00000
122	12 PHE		-37.00216		98.93077 A	1 12	0.00000
123	13 TYR		-35.16677	-20.22034	99.15293 A		0.00000
124	13 TYR	H	-34.71888	-21.10886	99.03348 £	1 13	0.00000

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125	13 TY	R CA	4.4465	0 -19.0289	1 99.5790	R A1	13	0.00000
126	13 TY			4 -19.3163			13	0.00000
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128				2 -17.5270			13	0.00000
129				6 -16.3571		-	13	0.00000
130 131				8 -17.4068 5 -16.2333			13 13	0.00000
132	•			8 -15.7159			13	0.00000
133	•			1 -14.5893			13	0.00000
134			-33.4734				13	0.00000
135				2 -18.5254		5 3 1	13	0.00000
136	13 TYP		-32.5945			5 A1	13	0.00000
137	14 LEU	-		0 -17.25550			14	0.00000
138	14 LEU			8 -16.68300			14	0.00000
139	14 LEU			8 -16.70428			14	0.00000
140 141	14 LEU	_		4 -16.21044 9 -17.29674			14 14	0.00000
142	14 LEU			3 -16.92524			14	0.00000
143	14 LEU			-17.54798			14	0.00000
144	14 LEU			7 -15.56813			14	0.00000
145	14 LEU	0	-32.31049	-14.71998	98.40126	λ1	14	0.00000
146	15 ASN	N	-30.73908	-15.58168	97.02166	Al	15	0.00000
247	15 ASN	н		-16.38705			15	0.00000
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149 150	15 ASN	CB		-14.93729			15	0.00000
151	15 ASN 15 ASN	CG OD 1		-14.18957 -13.98878	98.46652 99.50693		15 15	0.00000 0.00000
152	15 ASN	ND2		-13.74712	98.26721		15	0.00000
153	15 ASN	HD21		-13.93409	97.42633		15	0.00000
154	15 ASN	HD22	• .	-13.18778	98.98020	Al	15	0.00000
155	15 ASN	C		-13.70870	95.65633	Al	15	0.00000
156	15 ASN	0		-14.17939	94.92357		15	0.00000
157	16 PRC	N		-12.64312	95.25454		16	0.00000
158	16 PRO	CD	•	-12.14147	.93.88390		16	0.00000
159 160	16 PRO 16 PRO	CA CB		-11.89943 -11.32145	96.02164 94.90180		16 16	0.00000
161	16 PRO	CG	-28.40713	-11.06430	93.72211		16	0.00000
162	16 PRO	c		-10.84560	96.98393		16	0.00000
163	16 PRO	ō		-10.81095	98.14625		16	0.0000
164	17 ASP	N	-29.73099	9.96981	96.45979	A1	17	0.0000
165	17 ASP	н.		-10.06794	95.52595		17	0.00000
166	17 ASP	CA	-30.07647	-8.75629		λ1	17	0.00000
167 168	17 ASP 17 ASP	CB CG	-30.80318 -30.22601	-7.83976 -6.44601	96.20071 96.27578	Al 11	17 17	0.00000 0.00000
169	17 ASP	OD1	-29.42577	-6.10216	95.40955		17	0.00000
170	17 ASP	OD2	-30.58500	-5.71147	97.19272		17	0.00000
171	17 ASP	С	-30.91226	-8.96778	98.44177	Al	17	0.0000
172	17 ASP	Ο.	-3Q.52677	-8.65960	99.56331		17	0.00000
173	18 GIN	N.	-32.11780	-9.49744	90.20428		18	0.00000
174 175	18 GLN		-32.36542	-9.81984	97.29256		18	0.00000
176	18 GLN 18 GLN		-33.10696 -34.05728	-9.57864 -8.37464	99.27949		18 18 .	0.00000
177	18 GLN		-33.36307	-8.37464 -7.07517	99.58476		18	0.00000
178	18 GLN	-	-33.97880	-5.85158	98.94046		10	0.00000
179	18 GLN		-35.13776	-5.50918	99.12725		18	0.00000
180	18 GLN		-33.14378	-5.16835	98.16624		18	0.0000
181	18 GLN	H521	-32.19859	5.47544	97.99847		18	0.0000
182	18 GLN		-33.43475	-4.32502	97.72302		1 8	0.00000
	18 GLN		-33.83924		99.26964		18	0.00000
184 185	18 GLN		-33.74750 ·		98.32391		18	0.00000 0.00000
186	19 SER 19 SER			-11.16924 : -10.43617 :			19 19	0.00000
197	19 SER			-12.50616			19	0.00000
53	19 SER			-12.36114			19	0.00000
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,18	9 19 SER	OG 34.37064 -11.87727 103.10841 A1 19	0.00000
19		HG 33.E3984 -12.63499 103.40254 A1 19	0.00000
] ¢		C -35.97640 -13.29776 100.02197 A1 19	0.00000
15		0 -36.49518 -12.86400 99.00144 21 19	0.00000
19.		N -36.24917 -14.51787 100.53004 A1 20	0.00000
19		H -35.87834 -14.79673 101.41380 A1 20	0.00000
195		CA -37.08223 -15.47533 99.79192 A1 20	0.00000
190 197		-36.39897 -15.89747 100.44590 Al 20	0.00000
198		-39.10838 -15.10421 101.05171 A1 20	0.00000
199		-38.72023 -17.16722 100.25477 A1 21	0.00000
200		-38.03254 -17.84573 99.94375 Al 21	0.00000
201		TA -40.11538 -17.64629 100.27114 A1 21 TB -40.54163 -17.64420 98.79559 A1 21	0.00000
202			0.00000
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205		E1 -43.22010 -17.59671 96.43390 A1 21 E2 -41.23758 -18.43207 96.25678 A1 21	0.00000 0.00000
206	21 GLU C		0.00000
207	21 GLU 0		0.00000
208	22 PHE N	· · · -41.54286 -19.38051 101.22645 A1 22	0.00000
209	22 PHE H		0.00000
210	22 PHE C	A -41.84777 -20.63223 101.93252 x1 22	0.00000
211	22 PHE C		0.00000
212	22 PHE CO		0.00000
213	22 PHE CI		0.00000
214	22 PHE C		0.00000
215	22 PHE CI		0.00000
216	22 PHE CE		0.00000
217	22 PHE CZ		0.0000
218	22 PHE C	-43.25845 -21.11908 101.58028 A1 22	0.0000
219	22 PHE 0	-44.19436 -20.32985 101.55111 A1 22	0.0000
220 221	23 MET N	-43.39639 -22.43138 101.32298 A1 23	0.00000
222	23 MET H 23 MET CA	-42.59132 -23.03279 101.30500 A1 23	0.00000
223	23 MET CA 23 MET CB		0.00000
224	23 MET CG	• <u>• • • • • • • • • • • • • • • • • • </u>	0.00000
225	23 MET SD	-46.67153 -23.10231 97.33272 k1 23	0.00000 0.00000
226	23 MET CE	-47.47592 -24.67640 96.98955 A1 23	0.00000
227	23 MET C	-44.60710 -24.56281 101.21411 A1 23	0.00000
228	23 MET O	-43.70841 -25.13566 100.62115 A1 23	0.00000
229	24 PHE N	-45.43241 -25.30285 101.99622 A1 24	0.00000
230	24 PHE H	-45.17063 -26.26850 102.03772 A1 24	0.00000
231	24 PHE CA	-46.72021 -25.11140 102.69060 A1 24	0.00000
232	24 PHE CB	-47.06193 -23.70552 103.21139 X1 24	0.0000
233	24 PHE CG	-46.27878 -23.27691 104.43188 A1 24	0.00000
234 235	24 PHE CD1		0.0000
236	24 PHE CD2		0.00000
237	24 PHE CE1 24 PHE CE2		0.00000
238	24 PHE CZ		0.00000
239	24 PHE C	-43.89594 -22.39162 106.72434 A1 24 -47.88569 -25.61775 101.85776 A1 24	0.00000
240	24 PHE 0	-47.88569 -25.61775 101.85776 A1 24 -48.73152 -24.87778 101.36424 A1 24	0.00000
241	25 ASP N	-47.89055 -26.94593 101.74049 A1 25	0.00000 0.00000
242	25 ASP H	-47.22437 -27.52403 102.20704 A1 25	0.00000
243	25 ASP CA	-48.86501 -27.64762 100.90165 A1 25	0.00000
244	25 ASP CB	-48.21052 -27.81535 99.52046 A1 25	0.00000
245	25 ASP CG	-49.19634 -20.30012 98.47932 A1 25	0.00000
246	25 ASP OD1	-49.75750 -27.46672 97.77589 A1 25	0.0000
	25 ASP 002	-49.39656 -29.50745 98.38197 A1 25	0.00000
	25 ASP C	-49.18430 -28.98699 101.56183 Al 25	0.00000
	25 ASP O	-40.57317 -29.35714 102.56241 A1 25	0.00000
	26 PHZ N	-50.15819 -29.70261 101.00377 A1 26	0.00000
	26 PHE H 26 PHE CA	-50.60050 -29.38662 100.15349 A1 26	0.0000
	26 PHE CA	-50.56030 -30.99701 101.54898 A1 26	0.00000
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25		B 51.42376 -30.83090 102.82009 A1	26 0.000	
25		G -52.60356 -29.92038 102.57724 A1	26 0.000	
25 25		D1 -52.50398 -28.55674 102.91777 A1	26 0.0000	
25	`	D2 -53.70380 -30.42265 101.99011 A1	26 0.0000 26 0.0000	
25		E2 -54.86474 -29.55076 101.74257 A1	26 0.0000 26 0.0000	
25	9 26 PHE C	Z -54.76149 -28.18547 102.07912 A1	26 0.0000	
26		-51.29340 -31.82500 100.51070 A1	26 0.0000	0
263			26 0.0000	
263 263	-		27 0.0000	
264			27 0.0000 27 0.0000	_
265	_		27 0.0000	
266			27 0.0000	-
267		01 -52.95640 -36.23164 101.40272 A1	27 0.0000	
268 269		02 -54.89062 -36.19966 100.47314 A1 -51.70102 -34.05698 98.54820 A1	27 0.0000 27 0.0000	
270		-52.51335 -34.10386 97.63254 A1	27 0.0000 27 0.0000	
271	28 GLY N	-50.39045 -33.92246 98.35075 A1	28 0.0000	
272		49.78144 -33.80079 99.13160 A1	28 0.0000	
273	28 GLY C		28 0.0000	
274 275	28 GLY C 28 GLY O	-50.00090 -32.51821 96.28701 A1	28 0.0000	
276	28 GLY O 29 ASP N	-A9.04304;-32.03342 95.69731 A1 -51.21573 -31.96145 96.33541 A1	28 0.0000 29 0.0000	
277	29 ASP H	-51.93955 -32.39453 96.87595 A1	29 0.0000 29 0.0000	
278	29 ASP CA		29 0.0000	
279	29 ASP CB		29 0.0000	
280	29 ASP CG	-52.46245 -30.54960 93.07981 A1	29 0.0000	
281 282	29 ASP OD 29 ASP OD		29 0.0000	
283	29 ASP C	2 -53.56233 -30.10545 92.75048 A1 -52.21461 -29.64631 96.07233 A1	29 0.00000 29 0.00000	
284	29 ASP 0	-52.39419 -28.59407 95.46732 A1	29 0.00000	
285	30 GLU N	-52.65130 -29.81701 97.32405 A1	30 0.00000	
286	30 GLU H	-52.34793 -30.57647 97.90251 A1	30 0.00000	
287 288	30 GLU CA	-53.43516 -28.70871 97.86160 A1	30 0.00000	
289	30 GLU CB	-54.71008 -29.18152 98.54724 A1 -55.84100 -29.71506 97.67444 A1	30 0.00000 30 0.00000	
290	30 GLU CD	-57.13279 -29.50058 98.43828 A1	30 0.00000	
291	30 GLU OE1		30 0.00000	
292	30 GLU 052		30 0.00000	
293 294	30 GFA C	-52.69655 -27.80995 98.82987 A1 -52.13191 -28.20198 99.84476 A1	30 0.00000	
295	31 ILE N	-52.13191 -28.20198 99.84476 A1 -52.76110 -26.53268 98.46810 A1	30 0.00000 31 0.00000	
296	31 ILE H	-53.28532 -26.29454 97.65353 A1	31 0.00000	
297	31 ILE CA	-51.98442 -25.52556 99.18511 A1	31 0.00000	
298	31 ILE CB	-51.81933 -24.32331 98.23317 A1	31 0.00000	
299 300	31 ILE CG2	-53.16329 -23.65625 97.91607 A1	31 0.00000	
301	31 ILE CD	-50.75168 -23.33810 98.71478 A1 -50.41981 -22.27940 97.66271 A1	31 0.00000 31 0.00000	
302	31 ILE C	-52.51316 -25.12446 100.56211 A1	31 0.00000	
303	31 ILE O	-53.70233 -24.97567 100.82175 A1	31 0.00000	
304	32 PHE N	-51.54696 -24.95576 101.46562 A1	32 . 0.00000	
305 306	32 PHE H		32 0.00000	
305	32 PHE CA	• •	32 0.00000 32 0.00000	
308	32 PHE CG		32 0.00000	
309	32 PHE CD1	-50.23927 -27.18148 105.15120 A1	32 0.00000	
310	32 PHE CD2		32 0.00000	
311 312	32 PHE CE1 32 PHE CE2		32 0.00000	
313	32 PHE CZ	•	32 0.00000 32 0.00000	
314	32 PHE C		32 0.00000	
315	32 PHE 0	-52.76581 -22.45086 103.62205 A1	32 0.00000	
316	33 HIS N	-50.69098 -22.50451 102.70013 A1	33 0.00000	

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31			33 0.00000
31		A0.22576 -21.21727 103.23242 Al	33 0.00000
31° 320			33 0.00000
32		G -49.22984 -20.32142 105.45366 A1 D148.02826 -19.72793 105.37205 A1	33 0.00000 33 0.00000
322		D1 -47.28999 -19.92757 104.75923 A1	33 0.00000
323		02 -49.97214 -19.68828 106.45019 A1	33 0.00000
324		22 -49.20118 -18.70682 106.97086 A1	33 0.00000
325			33 0.00000
326		-49.11511 -20.78472 102.29629 A1 -48.59725 -21.61206 101.55845 A1	33 0.00000
327 328		-48.74537 -19.50449 102.33695 A1	33 0.00000 34 0.00000
329		-49.19429 -18.80802 102.90183 A1	34 0.00000
330			34 0.00000
331	34 VAL CE	-47.81303 -18.42567 100.28392 A1	34 0.00000
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335	34 VAL 0	-47.41849 -17.54298 103.42874 A1	34 0.00000 34 0.00000
336	35 ASP N-		35 0.00000
337	35 ASP H	-45.03039 -18.68906 101.76710 A1	35 0.00000
338	35 ASP CA	-44.68799 -16.93300 102.93355 A1	35 0.00000
.339 340	35 ASP CB	-44.74876 -15.73492 101.94639 k1	35 0.00000
341	35 ASP CG 35 ASP OD	-46.11850 -15.06364 101.88569 A1	35 0.00000
342	35 ASP OD:	· , ; · · · · · · · · · · · · · · · ·	35 0.00000 35 0.00000
343	35 ASP C	+44.90949 -16.57120 104.41152 11	35 0.00000
344	35 ASP O	-45.03304 -17.43137 105.27733 A1	35 0.00000
3 4 5	36 MET N	-44.91212 -15.26386 104.68945 X1	36 0.00000
346	36 MET H .	-44.97503 -14.58792 103.95356 A1	36 0.00000
347 348	36 MET CA 36 MET CB	-45.05621 -14.74244 106.04065 A1	36 0.00000
349	36 MET CB 36 MET CG.	-44.58443 -13.28473 106.01845 A1 -43.15122 -13.10955 105.50403 A1	36 0.00000 36 0.00000
350	36 MET SD	-42.96722 -11.71848 104.37016 Al	36 0.00000
351	36 MET CE	-43.60204 -10.40582 105.42531 A1	36 0.00000
352	36 MET C	-46.49207 -14.77022 106.53712 A1	36 0.00000
353 354	36 MET 0	-46.75425 -14.73983 107.73458 A1	36 0.00000
355	37 ALA K 37 ALA H	-47.43476 -14.78996 105.58618 A1 -47.21869 -14.87225 104.60519 A1	37 0.00000 37 0.00000
356	37 ALA CA	-48.80100 -14.57493 106.05297 A1	37 0.00000
3,57	37 ALA CB	-49.30059 -13.19637 105.61316 A1	37 0.00000
358	37 ALA C	-49.83252 -15.61256 105.65673 A1	37 0.00000
359	37 ALA O	-49.85254 -16.18787 104.57639 A1	37 0.00000
360 361	38 LYS H	-50.76933 -15.79391 106.59397 A1	38 0.00000
3 6 2	38 LYS H 38 LYS CA	-50.65212 -15.37866 107.49351 A1 -52.00981 -16.48765 106.23832 A1	38 0.00000 38 0.00000
363	38 LYS CB	-52.90628 -16.55867 107.48308 A1	38 0.00000
364	38 LYS CG	-52.41585 -17.40517 108.60236 A1	38 0.00000
365 366	38 LYS CD,	-53.40991 -18.62084 108.869B3 A1	38 0.00000
367	38 LYS NZ	-53.42547 -19.66155 107.75111 A1 -54.76503 -20.17861 107.56038 A1	38 0.00000 38 0.00000
368	38 LYS H21	-54.79226 -20.92377 106.83759 A1	38 0.00000 38 0.00000
369	38 LYS HZ2	-55.17879 -20.59510 108.44046 A1	38 0.00000
370	3B LYS H23	-55.42747 -19.41621 107.27224 A1	38 0.00000
371	38 LYS C	-52.74081 -15.73437 105.12989 A1	38 0.00000
372	38 LYS 0	-52.72521 -14.51209 105.10297 A1	38 0.00000
373 374	39 LYS N	-53.35457 -16.44351 104.17702 A1	39 0.00000
375	39 LYS H 39 LYS CA	-53.59774 -15.96113 103.33578 A1 -53.67982 -17.86931 104.22041 A1	39 0.00000 39 0.00000
376	39 LYS CB	-55.18971 -15.00506 104.55346 A1	39 0.00000
377	39 LYS CG	-56.24681 -17.84030 103.42417 A1	39 0.00006
378	38 TRS CD	-56.27039 -15.52580 102.62130 A1	39 0.00000
375	39 LYS CE	-56.64532 -15.69151 101.13213 A1	39 0.00000
350	39 LYS NZ	-55,66137 -17.54002 100.43802 A1	39 0.00000

./52	7 ಗುಟ್ಟು ಬಾ	Thu F	25 14:58 كت	:48 1993	7	
38		H21 55.95	645 -18.545	81 100.36313	λ1 3	0.00000
38:		H22 -55.39	863 -17.267	25 99.46103 .		0.00000
33					A1 39	0.00000
38	_				A1 39	0.0000
385				18 101.88936		0.00000
381 387				20 102.71885		0.00000
388				91 103.42697 <i> </i> 45 101.48663 <i> </i>		0.00000
389				43 100.44056		0.00000
390			03 -19.827			0.00000
391			27 -18.901			0.00000
392	40 GLU (DE1 -53.606	81 -17.683	72 98.89245 2	40	0.00000
393			46 -19.3556			0.00000
394				6 101.61501 2		0.00000
395				4 101.08479 A		0.0000
396 397	41 THR N			6 102.35754 A		0.00000
398				2 102.86321 A 4 102.30963 A		0.00000
399				0 102.36303 A		0.00000
400			42 -24.7540			0.00000
401	41 THR H		35 -25.5961			0.00000
402	41 THR C			3 103.43152 A		0.00000
403	41 THR C			0 103.42344 A		0.00000
404	41 THR O			6 104.56718 A		0.00000
405	12 VAL N			5 103.00455 A		0.00000
406 407	42 VAL H			2 102.09633 %		0.00000
408	42 VAL C			1 103.80812 A 6 102.98525 A		0.00000
409				6 102.98525 A 4 103.74041 A		0.00000 0.00000
410				8 101.63610 A		0.00000
411	42 VAL C			105.14170 A		0.00000
412	42 VAL O			106.20404 A		0.00000
413	43 TRP N	-60.2665	2 -24.07991	7 105.07466 A	43	0.00000
414	43 TRP H			7 104,20145 A		0.0000
415	43 TRP CA	-60.3517		3 106.34002 A		0.0000
416 417	43 TRP CE			106.17013 A		0.0000
418	43 TRP CO			104.73262 A		0.00000 0.00000
419	43 TRP CE			103.92127 A2		0.00000
420	43 TRP CE			104.12574 21		0.00000
421	43 TRP CD			103.93459 A1		0.00000
422	43 TRP NE		4 -27.31306			0.00000
423 424	43 TRP HE		0 -27.49046			0.00000
425	43 TRP CZ			101.58525 A1		0.00000
426	43 TRP CH			103.07694 A1 101.82597 A1		0.00000
427	43 TRP C			107.16564 A1		0.00000 0.00000
428	43 TRP O			108.02836 A1		0.00000
429	44 ARG N			106.89519 A1		0.00000
430	44 ARG H			106.07364 A1	44	0.00000
431	44 ARG CA			107.89144 %1	4 4	0.00000
432	44 ARG CB			107.26013 A1	44	0.0000
433 434	44 ARG CG			108.22026 A1	44	0.00000
435	44 ARG CD			107.47006 A1	44	0,00000
436	44 ARG HE			108.32672 A1 109.25139 A1	4 4 4 4	0.00000 0.00000
437	44 ARG CZ			107.86475 A1	4 4	0.00000
438	44 ARG NHI			108.56844 A1	4.4	0.00000
439	44 ARG HH1	1 -48.98076	-22.84912	108.24673 A1	44	0.0000
440	44 ARG HH1			109.51721 A1	44	0.00000
441	44 ARG NH2			106.69352 A1	44	0.00000
442	44 ARG HH2	1 -49.73591	-23.96008	106.35139 A1	44	0.00000
444	AA ARG C			106.14195 A1 108.38029 A1	14 44	0.00000 0.00000
- •		94.92131	-24.73637	100.33025 P.1	7 7	0.00000

	_							
್ಷ. /ಬಸ	KIN:2.C	ಸ್	Thu s	ద్ది 25 14:58	3:48 1993		8	
44			€.536	064 -21.392	01 108.885	43 ki	44	0.00000
441	_	א טב	-58.652	270 -21.556	86 108.185	48 A1	45	0.00000
447	_	_		46 -22.063			45	0.00000
449				579 -20.218 267 -19.465			45	0.00000
450			-59:150	56 -18.969	90 106.495	79 A1	45 45	0.00000
451	45 L1	במ כם:		34 -18.419			45	0.00000
452				31 -17.922			45	0.00000
453				18 -20.201			45	0.00000
454 455				62 -19.199 17 -21.360			45	0.00000
456				47 -22.147			4 6 4 6	0.00000
457		JU CÀ	-60.583	79 -21.473	17 111.864	81 A1	46	0.00000
458	46 GI		-61.478	17 -22.715	18 111.954	37 X1	46	0.00000
459 460	46 GL 46 GL		-60.808	81 -24.0230 64 -25.1183	85 111.517	72 Al	46	0.00000
461	46 GL			67 -25.4722			46 46	0.00000 0.00000
462	46 GL	U OE2	-62.248	94 -25.6063	6 112.522	80 Al	46	0.00000
463	46 GL	υc	-59.4869	98 -21.4932	3 112.9268	3 Al	46	0.00000
464 465	46 GL		-59.6095	59 -21.2222	6 114,1049	3 A1	46	0.00000
466	47 GL		-58.2788	30 -21.7931 12 -22.0665	.0 112.4422	IA DI	47	0.00000
.467	47 GL		-57 0815	55 -21.7386	4 113 2769	1A 01	47 47	0.00000
468	47 GL	у св	-55.8912	1 -22.0948	1 112.3905	9 21	47	0.00000
469	47 GLI		-55.9503	6 -23.5066	1 111.7897	7 31	47	0.00000
470 471	47 GLU 47 GLU		-55.5541				47	0.00000
472	47 GLU			6 -25.7440 9 -24.2787			47	·0.00000
473	47 GLU			7 -20.3815			47 47	0.00000 0.00000
474	47 GLU			0 -19.3574			47	0.00000
475	48 PHE		-56.8380	7 -20.4071	6 115.2471		48	0.00000
476 477	48 PHE 48 PHE			3 -21.2812			48	0.00000
478	48 PHE			7 -19.1364° 1 -19.34982			4 8 4 8	0.00000 0.00000
479	48 PHE			7 -18.4984			48	0.00000
480	48 PHE		-59.7112	2 -19.00320	117.4421	5 A1	48	0.00000
401 482	48 PHE 48 PHE	CD2 CE1		9 -17.19966			48	0.00000
483	48 PHE	CE2	-59.38940	7 -18.20670 7 -16.40286			48 48	0.00000 0.00000
484	48 PHE	CZ		-16.90878			48	0.00000
485	48 PHE	C	-55.39240	-18.46628	115.93777		48	0.00000
486	48 PHE	0		-19.07618			48	0.00000
487 488	49 GLY 49 GLY	N		-17.15361			49	0.00000
489	49 GLY	CV H	-54 23420	=16.72101 =16.30520	116.34575	Al Al	49 49	0.00000 0.00000
490	49 GLY			-16.86464			49	0.00000
491	49 GLY	0	-51.85815	-16.69886	116.10714	A1	4 9	0.00000
492 493	50 ARG			-17.56211			50	0.00000
494	50 ARG			-17.67284 -16.16800			50 50	0.00000
495	50 ARG			-18.92760			50	0.00000 0.00000
496	50 ARG	CG .	-51.10260	-19.32736	120.58254	Al	50	. 0.00000
497 498	50 ARG			-20.14774			50	0.00000
499	50 ARG			-20.44943			50	0.00000
500	50 ARG			-19.71957 -21.63184			50 50	0.00000 0.00000
501	50 ARG			-21.84466			50	0.00000
502	50 ARG	HH11 -	-49.05579	-22.71179	124.52612	A.1	50	0.00000
503 504	50 ARG			-21.13233			50	0.00000
505	50 ARG			-22.58629 -23.47414			50 50	0.00000 0.00000
506	50 ARG			-22.42795			50	0.00000
507	50 ARG	c -	51.06703	-19.09561	117.49074		50	0.00000
508	SO ARS	0 -	49.84240	-19.09133	117.41926	<i>λ</i> :	50	0.00000

. /ಏ೩	י באָבאַיַ.	೨೨	Thu Fab 25 14:58:4	8 1993	9	
50	9 51 p	HE N	. 84915 -19.87778		. 5i	0.00000
51	0 51 P	н эн	2.84564 -19.76657	116.76797 A	51	0.00000
511		HE C	-51.29477 -20.78796	115.73402 A	51	0.00000
512 513			-52.51672 -21.52503	115.15449 A1		0.00000
514			-52.25099 -22.49194 1 -51.87888 -23.82851			0.00000
515				114.2\$503 A1 112.69289 A1		0.00000
516			1 -51.72483 -24.74421	113.22969 A1	. 51	0.00000
517						0.00000
518			-51.94314 -24.31860			0.00000
519 520			-50.50268 -20.00696 -49.31281 -20.20801			0.00000
521			-51.20469 -19.03212		52	0.00000
522		LA H		114.33870 Al	52	0.00000
523	52 AI		-50.54896 -18.16486		52	0.00000
524	52 AI			112.62511 A1	52	0.00000
525 526	52 AI 52 AI		-49.28257 -17.48933 -48.27008 -17.39834	113.62687 A1	52	0.00000
527	53 SE		-49.35763 -17.04955 :		52 53	0.00000
528	53.55			115.38567 21	53	0.00000
529	53 SE	R CA	-48.18100 -16.47055	115.53815 A1	53	0.00000
530	53 SE		-48.58146 -16.02158 1	116.95383 A1	53	0.00000
531 532	53 SE 53 SE		-47.59593 -15.15541 1	117.52845 A1	53	0.00000
533	53 SE	_	-47.05156 -14.90357 1 -46.99433 -17.42737 1	118.42105 A1	53 53	0.00000
531	53 SE		-45.89463 -17.11790 1	15.12568 A1	53	0.00000 0.00000
535	54 PH		-47.26082 -18.64220 1	16.08200 A1	54	0.00000
536	54 PH			16.41568 A1	54	0.00000
537 538	54 PHE 54 PHE		· ·	16.09999 A1	54	0.00000
539	54 PHS			16.63413 A1 18.13255 A1	5 4 5 4	0.00000
540	54 PHE		-48.11656 -21.50316 1	18.65539 AI	54 54	0.00000 0.00000
541	S4 PHE			19.00621 A1	54	0.00000
542	54 PHE			20.05266 A1	54	0.00000
543 544	54 PHE 54 PHE			20.40382 A1	54	0.00000
545	54 PHE		-47.30008 -21.09192 1: -45.57270 -19.90650 1:		54	0.00000
546	54 PHE			14.73758 A1 14.55246 A1	54 54	0.00000 0.00000
547	55 GLU	N	-46.46681 -20.08976 13		55	0.00000
5 4 8	55 GLU	-	-47.45338 -20.04226 13	3.95516 A1	55	0.00000
549	55 GLU	CA		12.42886 A1	55	0.00000
550 551	55 GLU	CB CG		1.54876 A1 2.19274 A1	55	0.00000
552	55 GLU	.CD	-49.04456 -22.36202 11		55 55	0.00000 0.00000
553	55 GLU	051	-49.15553 -23.53977 11	.0.96047 Al	55	0.00000
554	55 GLU	OE2	-49.80524 -21.50984 11	0.81474 A1	55	0.00000
555 556	55 GLU 55 GLU	c:	-45-19354 -19.30188 11	1.76318 A1	55	0.00000
557	56 ALA	N .	-44.12678 -19.50387 11 -45.73650 -18.08673 11	1.19343 A1	55	0.00000
558	56 ALA	E	-46.62336 -17.96765 11	1.69532 A1 2 34890 A1	56 56	0.00000 0.00000
559	56 ALA	CA	-45.00414 -16.92733 11	1.38640 A1	56	0.00000
560	56 YTY	CB	-45.80074 -15.63871 1i	1.59969 AL	56	0.0000
561 562	56 ALA	C	-43.63772 -16.77849 11	2.02791 A1	56	0.00000
563	56 ALA 57 GLN	О И	-42.62065 -16.60634 11 -43.63088 -16.80634 11	1.36878 A1	56	0.00000
564	57 GLN	н	-43.63088 -16.91454 11: -44.48826 -17.03983 11:	3.33327 Al 1. 05838.5	57 57	0.00000 0.00000
565	57 GLN	Cλ	-42.35063 -16.88200 11	4.06847 Al	57	0.00000
566	57 GLN	СЭ	-42.61967 -17.03049 11:	5.56930 Al	57	0.00000
567 566	57 GLN	CG	-{1.40659 -16.75355 116	6.45895 Al	57	0.0000
569	57 GLN 57 GLN	CD	41.80070 -16.94589 117	7.90800 21	57	0.00000
	57 GLN		-42.29879 -16.06206 118 -41.55425 -16.16011 118		57 57	0.00000 0.00000
571	57 GLN		41.14709 -18.27176 117		57 57	0.00000
572	57 GIA		41.78115 -25.36728 119		57	0.00000

FIG. 10

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573	57 CLN C	1.36213 -17.94055 113.59025 A1	57 0.00000
574		-40.18971 -17.68051 113.34773 A1	57 0.00000
373	58 GLY N	-41.88825 -19.15742 113.41319 h1	0.00000 8 ق
576	58 GLY H	-42.84771 -19.33514 113.64847 A1	58 0.00000
577		-41.05379 -20.23393 112.87374 A1	58 0.00000
578		-40.45534 -19.92701 111.50957 A1	58 0.00000
579		-39.25324 -20.01021 111.27849 A1	58 0.00000
580		-41.34654 -19.52227 110.60018 A1	59 0.00000
581	59 ALA H	-42.32072 -19.46762 110.83485 A1	59 0.00000
582	59 ALA CA	-40.88666 -19.14537 109.26181 A1	59 0.00000
5B3	59 ALA CB	-42.07124 -18.74086 108.38101 A1	59 0.00000
584	59 ALA C	-39.86090 -18.02257 109.26073 A1	59 0.00000
585	59 ALA O	-38.85187 -18.04905 108.56603 A1	59 0.00000
586 587	60 LEU H	-40.12979 -17.02925 110.11113 A1 -40.96678 -17.03715 110.66448 A1	60 0.00000
588	60 LEU CA	-39.17026 -15.93079 110.21454 A1	60 0.00000
589	50 LEU CB	-39.82974 -14.72108 110.88234 A1	60 0.00000
590	60 LEU CG	-41.00342 -14.16448 110.06267 A1	60 0.00000
591	60 LEU CD		60 0.00000
592	60 LEU CD2	40.54968 -13.69403·108.67852 A1	60 0.00000
593	60 LEU C	-37.86300 -16.28607 110.90549 A1	60 0.00000
594	60 LEU O	-36.81366 -15.71151 110.64266 A1	60 0.00000
595	61 ALA N	-37.92548 -17.30628 111.76650 A1	61 0.00000
596	61 ALA H	-38.80416 -17.70206 112.04737 A1	61 0.00000
597	61 ALA CA	-36.66060 -17.86080 112.25036 A1	61 0.00000
598	61 ALA CB	-36.90091 -18.87147 113.37402 A1	61 0.00000
559	61 ALA C	-35.86652 -18.52303 111.13575 A1	61 0.00000
600	61 ALA 0	-34.67753 -16.28483 110.93578 A1	61 0.00000
601 602	62 A9N N 62 ASN H	-36.59182 -19.33811 110.35468 A1	62 0.00000
603	62 ASN K 62 ASN CA	-37.55651 -19.52444 110.56450 Al	62 0.00000 62 0.00000
604	62 ASN CA	-35.93048 -19.97053 109.20954 A1 -36.90608 -20.83397 108.41185 A1	62 0.00000 62 0.00000
605	62 ASN CG	-36.14296 -21.97501 107.76767 A1	62 0.00000
606	62 ASN 0D1	-35.90083 -23.00051 108.38915 A1	62 0.00000
607	62 ASN ND2	-35.81296 -21.80385 106.49294 A1	62 0.00000
608	62 ASN HD2.	•	62 0.00000
609	62 ASN HD2	•	62 0.00000
610	62 ASN C	-35.27272 -18.97317 108.27635 A1	62 0.00000
611	62 YZN 0	-34.08977 -19.05772 107.98073 A1	62 0.00000
612	eg ite k	-36.07385 -17.96130 107.91224 A1	63 0.00000
613	63 ILE H	-37.03805 -17.97906 108.18770 A1	63 0.00000
614	63 ILE CA	-35.60960 -16.86395 107.05550 A1	63 0.00000
615 616	63 ILE CB	-36.79680 -15.88630 106.84927 A1	63 0.00000
617	63 ILE CG2	-36.71800 -14.58751 107.66300 X1 -37.00443 -15.60068 105.36486 X1	63 0.00000
618	63 ILE CD	-36.27181 -14.79756 105.06416 $\lambda1$	63 0.00000
619	63 ILE C	-34.32421 -16.14412 107.48562 A1	63 0.00000
620	63 ILE 0	-33.67028 -15.43835 106.72047 Al	63 0.00000
621	64 ATA N	-33.97867 -16.34078 108.76481 A1	64 0.00000
622	64 ALA H	-34.55914 -16.88095 109.37800 A1	64 0.00000
623	64 ALA CA	-32.68252 -15.86370 109.23001 Al	64 0.00000
624	64 NIA C3	-32.78414 -15.37252 110.67448 X1	64 0.00000
625	64 ALA C	-31.59324 -16.91956 109.13883 Al	64 0.00000
626	64 AIA O	-30.45701 -16.65389 108.75412 A1	64 0.00000
627	65 VAL N		65 0.00000
628	65 VAL H		65 0.00000
625	65 VAL CA		65 0.00000
630 631	65 VAL CB		65 0.00000
632	65 VAL CG1 65 VAL CG2	* · · · · · · · · · · · · · · · ·	65 0.00000 65 0.00000
633	65 VAL C		65 0.00000
634	65 VAL 0		65 0.00000
635	65 ASP N		65 0.00000
636	56 ASP H		66 0.00000

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63	ر مد 66 · 7	P CA	11.048	58 -10 937	16 105.7679		76.6	0.0000
638		P CB	- 2.310	36 -20 249	66 104.9300	2 71	66	0.00000
533		P CG			19 105.0820		66	0.00000
640					14 105.2072		66	0.00000
541					68 105.0812		66	0.00000
642					49 105.1556		66	0.00000
643		PÖ			72 104.5070		66	0.00000
644					16 105.4801		67	0.00000
645	67 LYS	S H			55 105.8405		67	0.00000
646	67 LYS	S CA			04 105.1932		67	0.00000
647	67 LYS				96 105.8710		67	0.00000
648	67 LYS	G CG			77 105.5685		67	0.00000
649					56 106.3816		67	0.00000
650	67 LYS	CE			45 106.14720		67	0.00000
651	67 LYS	NZ			91 107.01869		67	0.00000
652	67 LYS	HZl	-30.3326		98 106.85909		67	0.00000
653	67 LYS				61 106.80064		67	0.00000
654	67 LYS				04 108.01156		67	0.00000
655	67 LYS		-20.2811	7 -16.5809	93 105.64383	Al	67	0.00000
656	67 LYS				78 104.87661		67	0.00000
657	68 ALA	N			73 106.92795		68	0.00000
658	68 YIA	ъ.			5 107.53692		68	0.00000
659	68 ALA	CA	-267635	2 -17.1814	3 107.40958		68	0.00000
660	68 ALA	CB		7 -17.4484			68	0.00000
662	68 ALA	C			4 106.70618		68	0.00000
((3	68 ALA 69 ASN	0	-24.8098	9 -18.3329	7 106.37143		68	0.00000
664	69 ASN 69 ASN	N		7 -19.3747			69	0.00000
665	69 ASN	H		6 -19.3442			69	0.00000
666	69 ASN	CA CE		6 -20.5473			69	0.00000
667	69 ASN	CG		4 -21.7056 4 -22.2021			69	0.00000
668	69 ASN	OD1			5 107.12937 7 107.30600		69 69	0.00000
669	69 ASN	ND2		-21.9892			69	0.00000
670	69 ASN	HD21		-21.5311			69	0.00000
671	69 ASN				109.04652		69	0.00000
672	69 ASN	C			7 104.36379		69	0.00000
673	MZA 09	Ö			6 103.89106		69	0.00000
67.4	70 LEU	N			103.71664		70	0.00000
675	70 LEU	H			104.12686		70	0.00000
676	70 LEU	CA.	-25.93555	-18.80513	3 102.42930	Al	70	0.0000
677	70 LEU	CB	-26.70466	-17.57714	101.93156	A1	70	0.00000
678	70 LEU	CG	-28.07464	-17.87907	101.32608	A1	70	0.00000
679	70 LEU				101.23109		70	0.00000
680						X 1	70	0.00000
681	70 LEU	C	-24.47328	-18.42736	102.51389	Al	70	0.00000
682 683	70 LEU	0	-23.64160	-1B.86456	101.72791	A1	70	0.0000
684	71 GLU	N ·	-24.17065	-17.62592	103.54240	አ 1	71	0.0000
685	71 GLU 71 GLU				104.19342		71	0.00000
686	71 GLU				103.73624		71	0.00000
687		CG -	-22.88033	-15.23004	104.88750	^_ } }	71 71	0.00000 0.00000
688	• •				104.93364		71	0.00000
689					106.76970		71	0.00000
690					106.33423		71	0.00000
691					103.97868		71	0.00000
692	_				103.36820		71	0.00000
693					104.86364		72	0.00000
694							72	0.00000
695			21.56703		105.13649		72	0.00000
696					106.17923		72	0.00000
697	72 ILS (21.51796				72	0.0000
696		CG1 -	22.55172	-20.59465	107.45409	P. 2	72	0.0000
699					108.51186		72	0.00000
700	72 ILE (c -	21.22106	-21.35113	103.89490	- 1	72	0.00000

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70	72	ILE O	20.06679 -21.68162 103.64178 A1	72	0.00000
70	2 73 1	MET N	-22.24406 -21.67127 103.09738 Al	73	0.00000
70		H TI	-23.17806 -21.34412 103.27303 $\lambda 1$	73	0.00000
70		ati C	$-21.88577 -22.51146 101.95532 \lambda1$	73	0.00000
70		EI C	3 -23.02917 -23.44601 101.56487 A1	73	0.00000
70			-23.30745 -24.44882 102.69030 A1	73	0.00000
70 70				73	0.00000
70	-	_		73	0.00000
70.			-21.30840 -21.79430 100.75010 A1	73	0.00000
71		et o	-20.54740 -22.36938 99.97894 A1	73	0.0000
71		HR H	-21.60342 -20.49214 100.63510 A1 -22.28955 -20.04403 101.21510 A1	74	0.00000
713		HR CA		74	0.00000
. 714		HR CB		74 74	0.00000
715				74	0.00000 0.00000
716		HR HG		74	0.00000
717		IR CG		74	3.00000
718		HR C	-19.35170 -19.61359 100.02467 A1	74	0.00000
719		-	-18.48554 -19.49348 99.16750 A1	74	0.00000
720				75	0.00000
721	75 LY		-19.81875 -19.65220 102.02321 A1	75	0.00000
722	75 LY		-17.69833 -19.83955 101.77078 A1	75	0.0000
723 724	75 LY 75 LY		-17_6140819.57287 103.27797 A1	75	0.0000
725			-16.20208 -19.66153 103.86107 A1	75	0.00000
726	75 LY 75 LY		-16.21091 -19.57437 105.38567 A1	75	0.00000
727	75 LY		-14.81504 -19.74395 105.98417 A1	75	0.00000
728	75 LY.		-14.91179 -19.69891 107.44939 A1	75	0.00000
7.29	75 LY.		-13.96820 -19.83508 107.86470 A1 -15.29159 -18.77551 107.74094 A1	75	0.00000
730	75 LY		-15.54844 -20.45311 107.77639 A1	75	0.00000
731	75 LY		-17.14118 -21.21778 101.45102 A1	75 75	0.00000 0.00000
732	75 LYS	S 0	-16.11623 -21.36022 100.79660 A1	75	0.00000
733	76 ARC	מכ	-17.86151 -22.25466 101.90990 A1	76	0.00000
734	76 ARC	H	-18.69512 -22.09296 102.44436 A1	76	0.00000
735	76 ARC		-17.35520 -23.61422 101.67525 A1	76	0.0000
736	76 ARC		-18.33654 -24.68494 102.17145 A1	76	0.00000
737 738	76 ARG		-18.82789 -24.66219 103.62361 A1	76	0.00000
739	76 ARG		-19.55383 -25.98411 103.91795 A1	76	0.00000
740	76 ARG		-20.36155 -25.99642 105.14334 A1	76	0.00000
741	76 ARG		-21.34527 -25.85364 105.02467 A1 -19.85069 -26.28686 106.34915 A1	76	0.0000
742	76 ARG		-19.85069 -26.28686 106.34915 A1 -20.67426 -26.45770 107.38550 A1	76	C.00000
743	76 ARG	RELL	-20.32038 -26.64144 108.30551 A1	76 76	0.00000 0.00000
744	76 ARG	HH12	-21.67206 -26.42100 107.26619 A1	76	0.00000
745	76 ARG	NH2	-18.53304 -26.41209 106.51158 A1	76	0.00000
746	76 ARG	HH21	-18.12377 -26.62371 107.39854 A1	76	0.00000
747	76 ARG	HH22	-17.92942 -26.29542 105.72237 11	76	0.00000
748 749	76 ARG	С	-17.06072 -23.91761 100.20901 A1	76	0.0000
750	76 ARG	0	-16.02715 -24.45819 99.83552 A1	76	0.00000
751	77 SER	ĸ	-18.01091 -23.50462 99.36716 A1	77	0.0000
752	77 SER	H CA	· ·	77	0.00000
753	77 SER	CB		77	0.00000
754	77 SER	OG		77 77	0.00000
755	77 SER	HG	_ * * _	7 7 77	0.00000
756	77 SER	c		, , 77	0.00000 0.00000
757	77 SER			, , 77	0.00000
	78 ASN			, , 7	0.00000
	78 ASN	H	-16.44843 -21.84082 98.96113 A1	78	0.00000
760	78 ASN	CV	-15.74721 -20.53857 97.45264 A1 :	7 a	0.0000
	78 ASN	CB	-14.33390 -21.01377 97.09195 A1 1	7 E	0.0000
	75 ASN			7 0	0.0000
	78 ASN		-12.89966 -20.31632 98.89264 A1 3	7 E	0.00000
. 61	78 ASN	ND3	-13,552f4 -22,45457 98,85326 A1 T	8	0.00000

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7 €	5 78 ASN H		• • • • •			
76	/ mm	D21 4.084	12 -23.191	70 98.43678 A		0.00000
7.6		1022 -13.038	71 -22.653	62 99.68493 A		0.00000
76		-16.341	34 -19.721			0.00000
	20	-15.714	78 -19.429			0.00000
76		-17.596	12 -19.325	97 96.52354 A	L 79	0.00000
77			24 -19.557	71 97.38252 A		0.00000
77:		λ -18.2140	08 -18.489	18 95.49858 A	79	0.00000
77:	79 TYR C	B -19.6481	74 -18.9264			0.00000
77:	79 TYR C	G -19.7448	85 -20.3276			0.00000
774	79 TYR C	01 -20.6424	46 -21.2327			0.00000
775			7 -22.5418			0.00000
776			1 -20.7211			0.00000
777			4 -22.0333			
778						0.00000
779						0.00000
780	_:		5 -24.2343		79	0.00000
781			2 -24.3067		79	0.00000
	79 TYR C		5 -17.0126		79	0.00000
782	79 TYR 0		9 -16.5127		79	0.00000
783	80 THR N		7 -16.3323		80	0.00000
784	80 THR H		9 -16.7950		80	0.00000
785	80 THR CA	-17.0957	7 -14.9266	95.20513 A1	80	0.00000
766	80 THR CB	-15.8907	2 -14.8771	96.18098 Al	80	0.00000
787	80 THR OG		9.,-13.52440		80	0.00000
788	BO THR HG		0 -13.5015		80	0.00000
789	80 THR CG		9 -15.72265		80	0.00000
790	BO THR C		-14.32258	93.83840 21	80	0.00000
751	80 THR O		7 -14.97644			
792	81 PRO N		-13.08096		80	0.00000
793		-11.43243	713.08090	93.61451 A1	81	0.00000
794		-10.04687	-12.25452	94.50529 A1	81	0.00000
		-10.95963	3 -12.43774	92.32255 A1	81	0.00000
795	81 PRO CB	-18.08102	! -11.39225	92.28996 Al	81	0.00000
796	81 PRO CG	-18.22970	-10.94901		81	0.00000
797	81 PRO C		-11.80328	92.25040 A1	81	0.00000
798	81 PRO 0	-15.41926	-10.58936	92.16776 A1	81	U.00000
799	82 ILE N .			92.27154 A1	82	0.00000
800	82 ILE H	-14.69364	-13.66392	92.31920 A1	82	0.00000
801	82 ILE CA	-13.18946	-12.18130	92.13921 A1	82	0.00000
802	82 ILE CB	-12.60010	-11.87598	93.53959 A1	82	0.00000
803	82 ILE CG2		-13.13773	94.38674 A1	82	0.00000
804	82 ILE CG1	-11 31152	-11.05228	93.44331 A1		
805	82 ILE CD	-10.76554	-10.62709	94.80896 A1	82	0.00000
806	82 ILE C		-13.19907		82	0.00000
B 07		1 -11.36717		91.37376 A1	82	0.00000
808		2 -11.36/1/	-12.81/4/	90.75062 A1	82	0.00000
809		2 -12.72556	-14.3/446	91.38671 A1	82	0.00000
810		-17.53322	-0.31236	94.99084 B1	1	0.00000
811		-17.21994	0.44323	94.35235 Bl	1	0.00000
	83 GLY HT2	-16.86357	-1.12219	94.96444 Bl	1	0.00000
812	83 GLY HT3	-17.61098	0.01920	95.97150 B1	1	0.00000
813	83 GLY CA	-18.79853	0.91116	94.55151 B1	1	0.00000
814	83 GLY C	-18.52573	-2.38203	94.66351 B1	1	0.00000
815	83 CLY O	-17.35786	-2.70920	94.84086 B1	1	0.0000
816	84 ASP. N	-19.57260	-3.20239	94.59303 Bl	2	0.00000
817	84 ASP H	-20.49658	-2.89510	94.35600 B1	2	0.00000
818	84 ASP CA	-19.43900	-4.63200	94.86181 B1	2	0.00000
819	84 ASP CB	-19.44643	-4.83356	96.38475 B1	2	
820	84 ASP CG	-18.89301				0.00000
821	84 ASP OD1		-6.19619	96.73815 B1	2	0.00000
822		-17.69451	-6.29764	96.98109 B1	2	0.0000
	84 ASP OD2	-19.66566	-7.14958	96.75715 B1	2	0.0000
923 =24	84 ASP C	-20.62766	-5.31072	94.19848 B1	2	0.0000
824	84 ASP O	-21.46903	-4.61697	93.63479 B1	2	0.00000
925	85 THR N	-20.67796	-6.64606	94.24891 B1	3	0.00000
826	85 THR H	-20.04362	-7.15926	94.84032 B1	3	0.00000
£27	ES THR Ch	-21.75257	-7.39367	93.59154 21	3	0.00000
323	95 THR C5	-21.58903	-7.31950	92.05122 E1	3	0.00000
				•		

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82		THR OC	32.7	265	-7.8248	30 91.3996		3	0.00000
831	0 85 9	THR HO			-7.6015			3	0.00000
833			2 -20.32		-8.0207			3	0.00000
832			-21.74		-8.8324			3	0.00000
833	•		-20.76		-9.3037			3	0.00000
834			-22.86		-9.5277			4	0.00000
835			-23.62		-9.1376			4	0.00000
836			-22.93	360 -	10.8755	2 94.4483	0 B1	4	0.00000
837					10.7908		6 B1	4	0.00000
838					11.6536		2 81	4	0.0000
839					13.1029			4	0.00000
840			-22.12	580 -:	14.0391	97.2708	3 B1	4	0.00000
841	86 A		-21.93	083 -1	4.6883	6 96.5349		4	0.0000
842 843	86 A		-21.37	502 -1	4.0501	6 98.3861		4	0.00000
. 844	86 AI		-20.31	371 -1	4.85617	7 98.4475		4	0.00000
845	86 A			115 -1	4.87872	99.26299		4	0.00000
846	36 AI		2 -20.051	. y	5.47498	97.6902		4	0.00000
847	86 AF			11/ -1	3.26636	99.42563		4	0.00000
848				60 -1	3.26659) 100.26897) · 99.36831	B1	4	0.00000
849	86 AF		-23.735	77 -1	1 92065	27.2001		4	0.00000
850	86 AF		-24.852	44 - 1	1 52502			4	0.00000
851	87 PR		-23, 121					1	0.00000
852	87 PR		-21.732					5 5	0.00000
- 853	87 PR		-23.844	39 -1	4.02757	92.53087		5	0.00000
854	87 PR	ED 03	-22.785	28 -19	5.13066	92.39999		5	0.00000 0.00000
855	87 PR	.o cc	-21.434	6C -1	4.43376	92.52460	_	5	0.00000
856	87 PR	.o c	-25.103	90 -14	.54496	93.21975		5	0.00000
857	87 PR	0 0	-25.324	1 -14	.39544	94.41838		5	0.00000
858	88 AR	G N	-25.943	4 -15	.17061	92.39123	Bl	6	0.00000
859	98 AR	GH	-25.686	1 -15	.35588	91.44433		6	0.00000
8 60	BB ARG	•	-27.2321			92.90235	_	6	0.00000
861	80 ARC		-28.2907	e -15	.39805	91.81653	Bl	6	0.00000
862	BB ARC		-29.7086	3 -15	.55859	92.35844	B1	6	0.00000
B 63	88 ARC		-30.7915	0 -16	.91787	91.49236	B1	6	0.00000
864	88 ARG		-31.9274	1 -14	.57277	92.34607	Bl	6 .	0.0000
865 866	88 ARG		-31.7651	2 -14		93.33788		6	0.0000
867	OFA 88		-33.0888		.12165	91.86193		6	0.00000
868	88 ARG	•	-34.0656 -34.9544	3 -13 1 -13	. 21 /56	92.71322		6	0.00000
869	88 ARG		-33.9185	3 -13	.4017/	92.40144 93.69838		6	0.00000
870	88 ARG		-33.2662	3 -13	. 32333	90.54974	_	6 6	0.00000
871	88 ARG	KH21	-34,1273	7 -13	64013	90.16725		6	0.00000 0.00000
872	88 ARG	HH22	-32.5208	-14	19859	89.92214		6	0.00000
873	BB ARG	С	-27.2315	7 -17.	07404	93.36366		6	0.00000
874	88 ARG		-26.8959	-18.	00090	92.63574		6	0.00000
875	89 PHE	N	-27.6275	-17.	24057	94.62546	Bl	7	0.0000
876	89 PHE	H	-27.92346	-16.	46797	95.18173	B1	7	0.00000
877	89 PHE	CA	-27.64368	-18.	59443	95.17433		7	0.00000
878 879	89 PKE	CB	-26.56379	-18.		96.25427		7	0.0000
	89 PHE	CG	-25.20774	-18.	55805	95.61829 1		7	0.00000
880 881	89 PHE	CD1	-24.74915	-19.	48139	94.64858 E	-	7	0.00000
882	89 PHE	CD2	-24.42840	-17.	43023	95.96102	_	7 .	0.0000
883	89 PHE		-23.51045			94.00961 E	_	7	0.00000
	89 PHE		-23.18767			95.32384 E		7	0.00000
	89 PHE		-22.73800 -28.99632			94.34976 E		7	0.00000
	89 PHE		-26.99632 -29.85551			95.72084 E 95.94786 E		7	0.00000 0.00000
887	90 LED		-29.85352			95.87791 H		, B	0.00000
	90 LEU		26.36720	-20		95.76376 B			0.00000
889	90 LEU		30.48975			96.14113 B			C.00000
890	90 LEU		31.01265			94.64981 B			0.00000
	90 LEU	CG -	32.46079	-21.1	6692	94.43614 B			C.00000
292	90 LEU	CD1 -	32.81137	-21.9		93.19113 2			0.00000

FIG. 15

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893	90 LEU	CD2 3	.47498 -21.40	699 95.55222	B1 8	0.00000
894		c -30	.45467 -21.90	281 97.20252		0.00000
ذوة		0 -29	.74216 -22.85			0.00000
896			.28133 -21.71			. 0.00000
897		•	.78897 -20.85	· · · · _		0.00000
898	91 TRP	CA -31.	61477 -22.86			0.00000
899	91 TRP	CB -31.		789 100.54418	B1 9	0.00000
900	91 TRP		46050 -22.11	190 101.18157	B1 9	0.00000
901	91 TRP		22893 -22.79		B1 9	0.00000
902	91 TRP		30074 -21.96		B1 9	0.00000
903	91 TRP	CE3 -28.	77368 -24.003	327 100.60389	B1 9	0.00000
.904	91 TRP	CD1 -30.	26500 -20.964	27 101.96112		0.00000
905	91 TRP		99810 -20.870			0.00000
906	91 TRP		68566 -20.156			0.00000
907	91 TRP		97767 -22.405		B1 9	0.00000
908	91 TRP		44054 -24.404			0.00000
909	91 TRP		56124 -23.618			0.0000
910	91 TRP		96765 -23.395		_	0.00000
911	91 TRP		92486 -22.660			0.00000
912 913			02847 -24.715			
913	92 GLN 92 GLN		20362 -25.283	· · · ·		
915	92 GLN		33516 -25.347			
516	92 GLN		50105 -26.041			
917			43287 -25.057			0.00000
918		OE1 -32.9	12668 -25.793 99424 -25.888			0.00000
919	•		9268 -26.321			0.00000
920		HE21 -36.1				0.00000
921		HE22 -35.0		,		0.00000
922		C -34.3				0.00000
923				9 100.00960 1		0.00000
924		N -35.4				0.00000
925		H -36.2				0.00000
926		CA -35.5				0.0000
927	93 LEU (1178 -26.8150		11	0.00000
928	93 LEU (CG34.8	3577 -27.7076	6 103.99848 E	11	0.00000
929		CD1 -34.3	3031 -26.8382	4 105.14843 E	11	0.00000
930		DD2 -35.8	2812 -28.7416			0.00000
931		-36.9	•			0.00000
932			1692 -27.4757			0.0000
933	_		5045 -29.4171			0.00000
934	• •			6 101.79736 B		0.00000
935 936	•		1063 -30.1412			0.00000
937				0 100.13839 B		0.00000
938			1969 -32.4483	2 100.07544 B		0.00000
939)198 -32.4483)198 -33.4044			0.00000 0.00000
940			389 -33.9096			0.00000
941			798 -34.4199.			0.00000
942		· •	249 -33.1426			0.00000
943	* *	••	162 -34.5793	•		0.00000
944	94 LYS C	•		102.67642 B		0.00000
945	94 LYS 0		720 -31.9977			0.00000
946	95 PHE N		676 -30.85529			0.00000
947	95 PHE H		321 -30.05201			0.00000
948	95 PHE C			104.49053 B		0.0000
949	95 PHE C	39.13	251 -31.49226	105.86481 E		0.0000
950	95 PHE CO		104 -30.20819		1 13	0.0000
951	95 PHE CE		917 -28.9688 <i>6</i>	106.12996 B		0.00000
952	95 PHE CO	4	668 -30.26479	107.51173 B	1 13	0.0000
953	95 PHE CE			106.77985 BI		0.0000
954	95 PHE CE			108.16313 B1		0.0000
955	95 PHE CZ			107.79605 23		0.00000
956	95 PHE C	-41.220	005 -32.16799	104.57423 E	. 13	0.00000

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957	7 95 PH	0 3	0859	91 -31.3133	2 104.41691 1	31 13	-07-00000
958	96 GL				4 104.75822 1		0.00000
9.19				33 -34.0987			0.00000
960					9 104.40078 1		0.00000
961				5 -34.2537			0.00000 0.00000
962 963				0 -34.8796 5 -34.9249			0.00000
964				5 -34.1525			0.00000
965			-42.6796	2 -35.6855	9 100.23154		0.00000
966			-43.1312	9 -35.2539	3 105.17925 8	31 14	0.00000
967	96 GLU	0	-42.2751	0 -36.1139	5 105.37879 E	14	0.00000
9 6 B	97 CYS	N	-44.3962	1 -35.3443	1 105.62501 E	15	0.0000
969	97 CYS		-45.0528	1 -34.6056	6 105.44206 E	1 15	0.00000
970	97 CYS				7 106.25700 5		0.00000
971	97 CYS				0 107.76554 E		0.00000 0.00000
972 973	97 CYS 97 CYS	_			8 108.42288 E 6 105.55706 E		0.00000
974	97 CYS	0			6 104.88787 B		0.00000
975	98 HIS	N			1 105.71784 B		0.00000
976	98-HIS	H			2 106.26371 B		0.00000
977	98 HIS	CA			0 105.05785 B		0.00000
978	98 HIS	СВ	-46.0086	7 -40.6102	9 104.19437 B	1 16	0.00000
979	98 HIS	CG			7 102.97364 B		0.00000
.980	98 HIS	ND1			9 101.74183 B		0.00000
981	98 HIS	HD1			3 101.51181 B		0.00000
982 983	98 HIS	CD2 NZŽ			5 102.89430 B 1 101.58753 B		0.00000 0.00000
984	98 HIS	CE1			101.30733 B		0.00000
985	98 HIS	٥			106.05217 B		0.00000
986	98 HIS	ō			106.99225 B		0.00000
987	99 PHE	N			105.81218 9		0.00000
988	99 PHZ	H			105.00666 B		0.00000
989	99 PHE	CA			106.77662 B		0.00000
990	99 PHÉ	CB			107.59234 B		0.00000
991	99 PHE	CG			108.81497 B		0.00000
992	99 PHE	CD1			108.76543 B		0.00000 0.00000
993 994	99 PHE 99 PHE	CD2 CE1			109.87537 B		0.00000
995	99 PHE	CE2			111.08354 B		0.00000
996	99 PHE	CZ	-47.92137	-39.56658	111.02626 B	1 17	0.00000
997	99 PHE	Ë	-50.79242	-41.97404	106.19800 B	1 17	0.00000
998	33 bit£	ö		-41.71068	105.22101 B	1 17	0.00000
999	100 PEE	34	-50.86837	-43.12653	106.85844 B	1 18	0.00000
1000	100 PHE	Ή.	-50.32768	-43.27017	107.68944 B	18	0.00000 0.00000
1001	100 PHE	CA	-51.84/18	#44.10783 #45.57710	106.41132 B	1 18 1 18	0.00000
1003	100 PHE	CB CG	-51 08740	-46 34931	105.63001 B	18	0.00000
1004	100 PHE	CD1			105.55366 B		0.00000
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1008	100 PKE	CZ	-50.37408	-47.83859	103.35608 BI	. 18	0.0000
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1010	100 PHE 101 ASK	0			108.13319 B1 106.08672 B1		0.00000
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1013	101 ASK	CA			106.37273 B1		0.00000
1014	101 ASN	C3			105.52094 B1		0.00000
1015	101 ASN	CG	-57.86040	-44.83624	105.52911 21	19	0.00000
1016	101 ASH	OD 1	-58.43246	-44.06601	104.77523 21	19	0.00000
1017	101 ASH	KD2	-58.51327	-05.52184	206.41359 81	19	0.00000
1018	101 ASN	HD21	-58.04843	-46.21753	107.02623 21	19	0.00000
1015	101 ASH		-59.50666	~45.49785	106.46882 91	19 19	0.00000 0.00000
1020	101 %5::	С	-55.06277	-44.09262	107.83398 B1	1.7	0.0000

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1021 101 ASH 0 16.73737 -43.23997 108.39	9463 B1	19	0.00000
1022 102 GLY N -55.63460 -45.10941 108.40	6454 Bl	20	0.0000
102 102 GLY H -55.00370 -45.02188 108.03	1935 B1	20	0.0000
1024 102 GLY CA -55.97438 -45.32353 109.88	3157 B1	20	0.00000
1025 102 GLY C -55.02088 -44.59872 110.61	1953 B1	20	0.00000
1026 102 GLY 0 -54.46008 -45.17864 111.73	800 B1	20	0.00000
1027 103 THR N -54.84744 -43.29716 110.54	648 B1	21	0.00000
1028 103 THR H -55.35512 -42.90731 109.77	435 B1	21	0.00000
1029 103 THR CA -53.96549 -42.39542 111.30		21	0.00000
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1032 103 THR HG1 -54.81974 -40.09919 113.49		21	0.00000
1033 103 THR CG2 -54.86925 -42.72433 113.73	959 B1	21	0.00000
1034 103 THR C -52.56983 -42.92386 111.66	956 B1	21	0.00000
1035 103 THR O -51.96086 -42.59288 112.68	707 B1	21	0.00000
1036 104 GLU N -52.05837 -43.78433 110.78		22	0.00000
1037 104 GLU H -52.50836 -43.94021 109.90		22	0.00000
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1039 104 GLU CB -50.88647 -45.95191 111.05	395 B1	22	0.00000
1040 104 -GLU · CG ···51.13368 -46.62525 112.419	591 B1	22	0.00000
1041 104 GLU CD -50.07267 -46.21410 113.430	075 B1	22	0.00000
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1043 104 GLU 0E2 -48,89658 -46,13148 113,083	359 Bl	22	0.00000
1044 104 GLU C -49.56607 -43.91901 110.459	54 B1	22	0.00000
1045 104 GLU 0 -49.58628 -43.40196 109.348	94 B1	22	0.00000
1046 105 ARG N -48.46470 -44.04578 111.193	45 B1	23	0.00000
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1049 105 ARG CB -46.43619 -43.14957 112.091	R4 B1	23	0.00000
1050 105 ARG CG -45.11167 -42.37595 112.052	20 21	23	0.00000
1051 105 ARG CD -43.89570 -43.22016 111.650		23	0.00000
1052 105 ARG NE -42.65640 -42.48704 111.894	00 BJ	23	0.00000
1053 105 ARG HE -42.30777 -42.45915 112.839	43 B1	23	
1054 105 ARG CZ -42.05554 -41.77685 110.938	71 D1	23	0.00000
1055 105 ARG NH1 -41.02352 -41.00938 111.251		23	0.00000
1056 105 ARG HH11 -40.51476 -40.49136 110.568	03 B1		0.00000
1057 105 ARG RH12 -40.73208 -40.92747 112.219		23	0.00000
1058 105 ARG NH2 -42.47661 -41.82718 109.682	19 BI	23	0.00000
1059 105 ARG HH21 -42.19810 -41.15340 109.003	00 DI	23	0.00000
		23	0.00000
12.57.05		23	0.00000
		23	0.00000
	34 B1	23	0.00000
19152000 45.51527 100.5002		24	0.00000
1064 106 VAL H -46.72693 -42.61352 108.3718 1065 106 VAL CA -45.53349 -44.24499 107.5835	7 21	24	0.00000
1066 106 VAL CB -46.27081 -44.39073 106.2429	2 DI	24	0.00000
1067 106 VAL CG1 -45.79579 -45.65575 105.5236	2 DT	24	0.00000
1068 106 VAL CG2 -47.77990 -44.41166 106.4292	0 B1	24	0.00000
1069 106 VAL C -44.14065 -43.66075 107.4155	2 51	24	0.00000
17, 11005 -43, 00075 107, 4155	C 23	24	0.00000
	6 81	24	0.00000
12.0000 100.2311	9 BI	25	0.00000
	1 BI	25 .	0.00000
		25	0.00000
12:35003 43:01317 104:7003		25	0.00000
-41.30320 -42.02819 103.3838		25	0.00000
12.00023 202.2100		25	0.00000
1077 107 ARG NE -43.76838 -42.93515 102.2314	7 Bl	25	0.00000
1078 107 ARG HE -44.37946 -42.14312 102.2828	4 B1	25	0.00000
1079 107 ARG CZ -44.26510 -44.17490 102.2094	5 Bl	25	0.0000
1080 107 ARG NH1 -45.56598 -44.32747 102.3590	5 E:	25	0.0000
1001 107 ARG HHI1 -46.00629 -45.19207 102.34569	2 51	2.5	0.00000
1082 107 ARG HH12 -45.19539 -43.55730 102.54613		25	0.00000
1083 107 ARG HH2 -43.47983 -45.23695 102.05799		25	0.00000
1084 107 ARG HH21 -43.86541 -45.15896 102.0545	? 21	25	0.00000

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1085	s 107)	LRC HH	122 2.494	04 -45.119	97 101.94496	B1 25	0.00000
108		URG C			46 106.32504		0.00000
1000	107 🖈	RG O			81 106.43645		0.00000
1088		EU N			02 106.39627		0.00000
1089		EU H	,		85 106.29320		.0.0000
1090		EU CA			44 106.53751		0.00000
1091					54 108.01248 39 108.35874		0.00000
1092		en cd			29 109.53808		0.00000
1094		EU CD			0 108.50151		0.00000
1095					13 105.85764		0.00000
1096		EU O			3 105.60154		0.00000
1097	_		-39.521	50 -37.4782	8 105.56773	B1 27	0.00000
1098					7 105.72467		0.00000
1099	109 L				9 104.90334		0.00000
1100	109 1				9 103.39009		0.00000
1101	109 L				5 102.82634 1 102.17066		0.00000
1103	109 L				3 101.91138		0.00000
1104	109-L				1-105.21268		0.00000
1105	109 LE				2 105.47943		0.00000
1106	110 GI		-36.7009	2 -35.2976	8 105.12126	B1 20	0.00000
1107	110 GI		•		6 105.00992		0.00000
1108	110 GI				5 105.08184		0.00000
1109	110 GI				7 106.49578		0.00000
1110	110 GI	•			0 106.60993		0.00000
1111	110 GL	•			8 106.30413 1 1 106.10521 1		0.00000 0.00000
1113	110 GL				106.36491		0.00000
1114	110 GL	•			5 104.15500		0.00000
1115	110 GL	•			7 104.11585		0.00000
1116	111 AR				3 103.37222 1		0.00000
1117	111 AR				103.37387		0.00000
1118	111 AR	•		•	3 102.46249 E		0.00000
1119	111 AR				101.20352		0.00000
1120 1121	111 AR				100.75305		0.00000 0.00000
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1123	111 AR	,*			100.22757 E		0.00000
1124	111 AR	S CZ		-36.32384			0.00000
1125	111 AR		-38.78116	-36.68971	99.16119 E		0.00000
1126	111 ARG	•	• •	-37.49888			0.00000
1127	111 ARG			-36.16699			0.00000
1128 1129	111 ARC				98.41518 B 97.94325 B		0.00000 0.00000
1130	111 ARC				98.36264 B		0.00000
1131	111 ARG				102.08170 B		0.00000
1132	111 ARG				101.76699 B		0.00000
1133	112 CYS				102.12414 B		0.00000
1134	112 CYS				102.37455 B		0.0000
1135	112 CYS				101.73826 B		0.00000
1136	112 CYS				102.91930 B		0.00000
1137 1138	112 CYS				102.57697 B		0.00000
1138	112 C13				100.54380 B 100.40406 B		0.00000 0.00000
	113 ILE				99.65078 B		0.00000
	113 ILE	н	-31.84193		99.86111 5		0.0000
1142	113 ILE	CA.			98.32464 B		0.00000
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	113 ILE	CG2	-33.00339		97.82323 B		0.00000
	113 ILE	CG1	-31.52431		95.89635 B		0.00000
	113 ILE	CD	-32.44047		94.93272 B		0.00000
	113 ILE 113 ILE	C O	-30.08576		97.96536 B 98.21317 B		0.00000 0.00000
0	-:3 :16	J	-30.75333	-49.23300	30.21311 13.		0.50000

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114	9 114 24	R N	28.870	10 -27.255	64 97.41665 E	1 32	0.00000
115				70 -28.101		1 32	0.00000
115				36 -25.982		1 32	0.00000
115				88 -26.019	33 97.71345 B		0.00000
115				27 -24.7210			0.00000
115	4 114 TY	R CD1	-26.6277	70 -23.4582			0.00000
215	5 114 TY	R CEI	-25.8488	6 -22.2808			0.00000
115	6 114 TY	R CD2	-24.6182	24 -24.8021			0.00000
115	7 114 TY	R CE2	-23.8384	11 -23.6265	1 97.34588 B	1 32	0.00000
1158	8 114 TY	R CZ	-24.4560	0 -22.3694	0 \$7.53154 B	1 32	0.00000
1159	9 114 TY	R OH	-23.6896	7 -21.2191	7 97.50979 B	1 32	0.00000
1160	114 TY	п нн	-22.7967	6 -21.4301	5 97.21984 B	1 32	0.00000
1161	114 TY	R C	-28.1472	3 -25.8321	5 95.61430 B	1 32	0.00000
1162	2 114 TY	R O	-27.6637	5 -26.7096			0.00000
1163	115 ASI	N N	-28.6682	3 -24.6999	5 95.12919 B	1 33	0.00000
1164		H F	-29.0309	2 -24.0239	1 95.77643 B	1 33	0.00000
1165			-28.6376	2 -24.4154	5 93.68566 B	1 33	0.00000
1166			-27.2704	9 -23.8438			0.00000
1167				9 -22.4311			0.00000
1168					494.85370 B		0.00000
1169	_			5 -21.4835			0.00000
1170			1 -28.0488				0.00000
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1172	115 ASN			B -25.5930			0.00000
1173	•			0 -25.9140			0.00000
1174	116 CTN			1 -26.23583			0.00000
1175	Jiè civ			0 -25.9261			0.00000
1176	116 CIN			5 -27.38897			0.00000
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1185	116 GLN	0		-29.77893	•		0.00000
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1193	117 GLQ	OE2		-29.45965		35	0.00000
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1195	117 GLU	0	-28.39032	-29.54087	95.93593 B1	35	0.0000
1196	178 erá	N		-31.70267	95.31151 B1	36	0.0000
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1200	118 GLU	CG	-30.04821		96.04454 Bl	36	0.00000
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1202	118 G;U		-31.31366		96.62004 B1	36	0.00000
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1204	118 GLU		-27.45968		97.59954 B1	36	0.00000
1205 1206	118 GLU		-26.30375		97.22004 B1	36	0.00000
1207	119 SER		-27.77719		98.81671 B1	37 37	0.00000 0.00000
1207	119 SER		-28.73032		99.10568 51	37 37	0.00000
1208	119 SER		-26.67523		99.74333 B1	37	0.00000
1210	119 SER 119 SER				100.25438 B1 100.96654 B1	37	0.00000
1210						37	0.00000
1211	119 SER 119 SER				101.30354 51 100.87899 B1	37	0.00005
2	119 227		ראינרס.ט∡-	-229659	299.01977 21		0.0000

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121	3 119 SE	n •	2256	F -32 0055	4 101 4222	37	0.00000
121	-		15.7230	4 -32.9055	4 101.03211 B 8 101.67561 B	-	0.00000
121						-	0.00000
121		-			8 101.48344 B	=	0.00000
121			-27.8267		2 102.80099 B	-	
121					4 104.13888 B		0.00000
121					8 104.29388 B		0.00000
1219					0 104.33500 B		0.00000
1220	120 VA	L C			5 102.89372 B		0.00000
1223	120 VA	L O	-30.19812	2 -32.8865	6 102.49711 B	38	0.00000
1222	2 121 AR	3 N	-29.34164	4 -34.7535	6 103.41944 B	39	0.00000
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1224		G CA			2 103.49447 B		0.00000
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					102.86363 B1		0.00000
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1230					5 101.13259 B1		
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1233		нн12			3 102.71808 B1	39	0.00000
1234	121 ARG			-39.84536		39	0.00000
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. 1236	121 ARG					39	0.00000
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1239	122 PHE	N	-32.04075	-36.24095	105.26404 B1	40	0.00000
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1241	122 PHE	CA			106.24769 B1	40	0.00000
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1243	122 PHS	CG			108.30418 B1	40	0.00000
					108.77664 B1	40	0.00000
1244	122 PHE	CD1				40	0.00000
1245	122 PHE	CD2			109.19094 B1		
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1250	122 PHE	0			104.65587 B1	40	0.00000
1251	123 ASP	N			106.15389 B1	41	0.00000
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1253	123 ASP	CA			105.58825 B1	41	0.00000
1254	123 ASP	CB	-30.85171	-41.81726	105.76445 B1	41	0.00000
1255	123 ASP	CG	-29.93161		104.59405 B1	41	0.00000
1256	123 ASP	OD1	-28.81173	-41.11310	104.81757 B1	41	0.00000
1257	123 ASP	OD2	-30.34905	-41.80191	103.45856 B1	41	0.0000
1258	123 ASP	С	-33.32362	-41.63618	106.09965 B1	41	0.00000
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1260	124 SER	N			105.41429 B1	42	0.00000
1261	124 5ER	н	-32.95329	-43.10971	104.76900 B1	42	0.0000
1262	124 BER				105.58880 Bl	42	0.00000
1263	124 SER				104.46149 B1	42	0.00000
1264	124 SER				104.37671 B1	42	0.00000
1265	124 SER	HG	-36 01080	-44 R7590	105.27413 B1	42	0.00000
1266	124 SER				106.93634 B1	42	0.00000
1267	124 SER	0	-32.21640	-44 A7777	107.22372 B1	42	0.00000
1268					107.74330 B1	43	0.00000
	125 ASP					43	0.00000
1269	125 ASP				107.50494 B1		0.00000
1270	125 ASP				109.10471 B1	43	
1271	125 ASP	CB	-32.96210	-45.24580	109.50620 Bl	43	0.00000
1272	125 ASP	CG	-31.82155	-44.23245	109.55779 E1	43	0.00000
1273	125 ASP				108.89075 E1	43	0.00000
1274	125 ASP				110.29614 E1	43	0.00000
1275	125 ASP				110.10428 21	43	0.00000
1276	125 ASP	0	-35.10540 -	-43.79441	111.22872 21	43	0.00000

. /5	אד"אביא	2.CI		<u> </u>	ab 25 14	:58:4	8 1993		21	
12	77 12	6 VAI	N	51	515 -42.	26751	109 647	90 B1	44	5 .50500
12		_		209. دد-	39 -42.	12308	108.704		4.4	0.00000
12		6 VAI			346 -41.				44	0.00000
121				-36.048	314 -40.	82462	110.975	92 B1	4.4	0.0000
121			_		64 -39.				4 4	0.00000
128		6 VAI			94 -41.	00281	109.824	81 B1	44	0.00000
128		6 AYT 9 AYT		-33.576	25 -41.	13739	111.649	61 B1	4.4	0.00000
128		7 GLY			19 -40.7 41 -41.6				44	0.00000
128					98 -42.0				45 45	0.00000 0.00000
128					44 -41.8				45	0.00000
128			c		94 -41.5				45	0.00000
128	9 127		ō		55 -41.2				45	0.00000
129		GLU	N	-29.811	43 -41.6	5704	110.295	45 B1	46	0.00000
129		-	H		53 -42.1				46	0.00000
129			Cλ		56 -41.1		109.731		46	0.00000
129 129			CB CG		97 -42.2		109.1170		46	0.00000
129			CD		16 -42.2 31 -43.2		109.598		46 46	0.00000
129		·GLU		24.4964	16 -42.A	8073	108.8763	12 B1	46	0.00000
1291		GLU	OE2		0 -44.4		109.0034		46	0.00000
1290	128		С		6 -39.9				46	0.00000
1299	128	GLU	0		4 -39.5		108.3572		46	0.00000
1300	129	TYR	N		5 -39.42		108.3893		47	0.00000
1301	- •	TYR	H		8 -39.83		108.6862		47	0.00000
1302		TYR	CA.	-27.5001			107.4643		47	0.00000
1303		TYR	CB	-26.6384	2 -37.15	326	108.0156	0 B1	47	0.00000
1304		TYR	CC	-27.3085					47	0.00000
1305		TYR TYR	CD1	-26.6718					47	0.00000
1307		TYR	CD2	-27.2556 -20.5282					47 47	0.00000
1308	129	TYR	CE2	-29.1123	7 -33.07 5 -34.88	063 1	09.8357	1 B1	47	0.00000
1309	129	TYR	CZ	-28.4732	7 -34.75	726 1	11.0886	0 Bl	47	0.00000
1310	129	IYR	OH	-29.0500	5 -33.98	459 1	12.0722	l Bl	47	0.00000
1311	129	TYR	нн	-29.7039					47	0.00000
1312	129	TYR	C	-26.8253	1 -38.64	384 1	06.1594	l Bl	47	0.00000
1313 1314		TYR ARG	0	-25.6669					47	0.00000
1315		ARG	n H	-27.55686 -28.51666	-30.38 -38.10	162 1 451 1	05.08583		48	0.00000
1316		ARG	CA	-26.07326			03.80227		4 B 4 B	0.00000 0.00000
1317		arg .	СВ	-27.85650			02.71647		48	0.00000
1318	130	arg .	CG	-27.21143	-39.00	526 1	01.34112		48	0.00000
1319		ARG	CD	-28.23975			00.29808	Bl	48	0.00000
1320 1321	130		NE	-27.66322	-39.38	333	98.95629		48	0.0000
1322	130 . 130 :		HE CZ	-26.82589			98.82833		48	0.00000
1323	130		NH1	-28.29934 -27.82365	-39.963	348 .	97.93202		48	0.00000
1324	130			-28.26738	-40 239	13 .	96.69917 95.90996		4 B 4 B	0.00000 0.00000
1325	130 1		HH12	-27.01064	-39.252	97	96.53955		48	0.00000
1326		ARG I	NH2	-29.39843	-40.687	30	8.14392		48	0.00000
1327	130 /		HH21	-29.90446	-41.104	92 9	97.39118		48	0.00000
1328 1329	130 3			-29.72741			9.07957	Bl	48 .	0.00000
1330	130 x	ARG (3	-26.28084			3.45986		4 B	0.0000
1331	131 A) •	-26.96293 -24.95816			3.07688		48	0.00000
332				-24.45069			3.58668		49 49	0.00000 0.00000
333	131 A			-24.28607		94 10	2.98902		49	0.00000
334	131 A			-23.06137			3.80801		49	0.00000
1335		LA C	:	-23.85084	-36.206	33 20	1.58658	81	49	0.00000
1336		LA C		-23.17892			1.36065		۷9	0.00000
.337 .338	132 V 132 V			-24.28495	-35.387	74 10	0.63154		50	0.00000
339	132 V			-24.78984 -24.05930			0.85567		50	0.00000 0.00000
340	132 V	AL C		-24.03530 -25.12578			9.24192 8.36560		50 50	0.00000
		_								

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134:	1 132 VA			05 -35.4568		u 1	20	_0. <i>57</i> 00 3 0
134				48 -35.4885			50	0.00000
2247	••			46 -35.5339			50	0.00000
1344	4 132 VA			98 -36.1925			50	0.00000
1345		R N		46 -34.5356	7 99.35299	Bl	51	0.00000
1346				36 -34.0427			51	0.00000
1347				97 -34.1385			51	0.00000
1348				79 -33.0996 L9 -32.5670			51 51	0.00000
1349				19 - 32.3670 $18 - 32.1013$			51	0.00000
1351				0 -31.9694			51	0.00000
1352				5 -33.5975			51	0.00000
1353				7 -33.3095			51	0.00000
1354				9 -33.4421			52	0.00000
1355				2 -33.6572			52	0.00000
1356				8 -32.8873			52	0.00000
1357 1356	134 GLU			2 -32.81284 9 -34.1707			52 52	0.00000
1359	134 GLU			7 -34.04410			52	0.00000
1360				1 -34.96158			52	0.00000
1361	134 GLU	OE2		1 -33.03519			52	0.00000
1362	134 GLU		-18.1860	1 -31.51310	101.45728	Bl	52	0.00000
1363	134 GLU	0		4.,-31.11916			52	0.00000
1364	135 LEU	N		1 -30.80674			53	0.00000
1365	135 LEU 135 LEU	н.		2 -31.13340			53	0.00000
1366 1367	135 LEU	CA CB	-19.3572 -20.2991	7 -29.57029 4 -29.06218			53 53	0.00000 0.00000
1368	135 LEU	CG		i -27.57937			53	0.00000
1369	135 LEU	CD1		-26.68772			53	0.00000
1370	135 LEU	CD2		-27.23870			53	0.00000
1371	135 LEU	C		-29.72915			5 3	0.0000
1372	135 LEV	0.		-28.93514			53	0.00000
1373	136 GLY	N :		2 -30.82221			54	0.00000
1374	136 GLY	н.		-31.49656			54	0.00000
1375	136 GLY	CA		-31.10000 -31.85099			54 54	0.00000 0.00000
1376 1377	136 GLY 136 GLY	C o		-31.82013			54	0.00000
1378	137 ARG	N		-32.53644			55	0.00000
1379	137 ARG	Н		-32.57514			55	0.00000
1380	137 ARG	CA		-33.32437			55	0.00000
1381	137 ARG	CB		-34.02762			55	0.00000
1382	137 ARG	CG	-18.13008				55	0.00000
1383	137 ARG	j. B		-36.15719			55 55	0.00000
1385	137 ARG 137 ARG	ne . He		-37.06640 -36.71980			55	0.00000
1386	137 ARG	CZ		-38.26612			55	0.00000
1387	137 ARG	NH1		-39.07643			55	0.00000
1388	137 ARG			-39.99255			55	0.00000
1389	137 ARG			-38.75541			55	0.00000
1390	137 ARG	NH2		-38.64369			55	0.00000
1391 1392	137 ARG 137 ARG			-39.53705			55	0.00000
1393	137 ARG	HH22		-38.01490 -32.65026			5 5 55 ·	0.00000 0.00000
1394	137 ARG	0'		-33.13638			55	0.00000
1395	138 PRO	N		-31.51331			56	0.00000
1396	138 PRO	CD		-30.78028			56	0.00000
1397	138 PRO	CA		-30.86885		11	56	0.00000
1398	138 PRO			-29.66775			56	0.0000
1399	138 PRO			-30.01427			56	0.00000
1400	138 PRO			-30.39451			56	0.00000
1401	138 PRO			-30.20105			56 57	0.00000 0.00000
1402 1403	139 ASP			-30.19487 -30.45258			57 57	0.00000
1404	139 ASE			-30.45258			57	C.00000
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. /22	יז־אביז יכ	æ	ಮಾ ಕಿಂಶ	25 14:58	:48 1993		23	
140	5 139 A	SP CB	3.29483	-29.227	54 106.807	20 81	5 7·	_0~0 ∞ 000
140	6 139 A	SP CG		-28.441	50 107.234	06 91	57	0.00000
1.0		SP OD	1 -24.33925	-27.259	55 107.536	98 Bl	57	0.00000
140		SP 00	2 -25.59590	-29.009	37 107.248	91 81	57	0.00000
140		SP C	23.29009	-30.906	51 108.706	46 Bl	57	0.00000
141		5P 0		-30.799			57	0.00000
141		_			50 108.049	75 B1	58	0.00000
141				-32.093			58	0.00000
141					3 108.701		58	0.00000
141					5 107.801		58	0.00000
141					1 110.046		58	0.00000
141					0 111.080		58	0.00000
1417		-			0 110.015		59	0.00000
1418	8 141 GI				2 109.142		59	0.00000
1419	141 GL				5 111.2627		59	0.00000
1420) 141 GL				4 110.9263		59	0.00000
1421	141 GL	U CG			5 109.9718		59	0.00000
1422	141 GL	U CD			0 109.4467		59	0.00000
1423	141 GL	U OE1	-17.21662				59	0.00000
1424		U - 0E2	·· -16.99658	-33.8737	5 -109.5028	8 B1	59	0.00000
1425			-21.20315	-32.7239	5 112.3669	6 B1	59	0.00000
1426			-21.35204				59	0.00000
1427			-21.39109				60	0.00000
1428		R H	-21.16858				60	0.00000
1429	142 TY	R CA	-21.91640				60	0.00000
1430	142 TY		-22.17510	-29.18770	111.9547	8 B1	60	0.00000
1431	142 TY	R CG	-22.15441	-27.86866	112.6890	2 Bl	60	0.00000
1432	142 JY	CD1	-20.91930				60	0.00000
1433	142 TYF		-20.89216	-25.94177	113.4963	3 B1	60	0.00000
1434	142 TYP		-23.36373				€0	0.00000
1435	142 TYP		-23.33600			B B1	60	0.00000
1436	142 TYP		-22.10013			4 Bl	60	0.00000
1437	142 TYR		-22.06472				60	0.00000
1438	142 TYR		-22.95958				60	0.00000
1439	142 TYR		-23.20365 ·				60	0.0000
1440	142 TYR	_	-23.33185				60	0.00000
1441	143 TRR		-24.16819 -				61	0.00000
1443	143 TRP		-24.01203 -				61	0.00000
1444	143 TRP 143 TRP		-25.46084 -				61	0.00000
1445	143 TRP 143 TRP	CB CG	-26.46502 -		112.10045		61	0.00000
1445	143 TRP	CD2	-26.82927 -				61	0.00000
1447	143 TRP	CE2	-27.59514 - -27.68725 -		110.26190		61 61	0.00000 0.00000
1448	143 TRP.	CE3	-28.23171 -				61	0.00000
1449	143 TRP	CDI	-26.49404 -				61	0.00000
1450	143 TRP	NEI	-26.99373 -				61	0.00000
1451	143 TRP	HE1	-26.86335 -				61	0.00000
1452	143 TRP	CZ2	-28.41151 -				61	0.00000
1453	143 TRP	CZ3	-28.94655 -				61	0.00000
1454	143 TRP	CH2	-29.03488 -				61	0.0000
1455	143 TRP	C.	-25.40824 -				61	0.00000
1456	143 TRP		-26.13451 -				61	0.00000
1457	144 ABN		-24.46546 -				62	0.00000
1458	144 ASN		-23.94027 -				62	0.00000
1459	144 ASN		-24.16067 -				62	0.00000
1460	144 ASN		-23.20850 -				62	0.00000
1451	144 ASN		-23.89541 -				62	0.00000
	144 ASN		-23.60155 -3				62	0.00000
	144 ASN	ND2	-24.72776 -3	37.64097	113.02355	81	62	0.00000
	144 ASN	HD21	-24.09338 -3	37.61926	113.99276	81	62	C. C0000
	144 ASN	HD22 -	-25.19584 -3	88.20406	112.34455	Bl	62	0.00000
	144 ASN		-23.49875 -3				62	0.00000
	144 ASN		-23.43003 -3				62	0.00000
458	145 SER	1: .	-22.99604 -3	3.45640	115.97355	31	63	0.00000

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1469 145 E		22.978	30 -32.80	^{C28} 115.2196	1 B1 6	000000000000000000000000000000000000000
1470 145 s		22.380	04 -33.14	521 117 258 <i>6</i>	S B1 6	0.00000
1472 145 S		-20.919	21 -32.73	54 117.0117	6 B1 6	0.00000
1473 145 S		-20.189	55 -32.04.	180 118.2451 35 118.9106	1 81 6	3 0.00000
1474 145 S		-23.110	27 -32.076	14 118.0616	ם ומכי	3 0.00000 3 0.00000
1475 145 61		-22.674	01 -31.682	87 119.1394	1 B1 6	
1476 146 GI		-24.221	80 -31.590	37 117.5125	2 Bl 6	
1477 146 GI		-24.581	18 -31.946	04 116.6477	4 Bl 6	
1478 146 GI 1479 146 GI		-24.900	25 -30.492	86 118.1952	2 Bl 6	
1479 146 GI 1480 146 GI				01 117.2389		
1481 146 GI				45 117.7288: 62 117.2523		
1482 146 GI		-27.808	37 - 27.536	09 118.0091	(B) 64	
1483 146 GL	N NE2	-27.0655	56 -27.979	48 115.9432	9 B1 64	
1484 146 GL		-26.3080	9 -28.203	49 115.33276	S B1 64	
1485 146 GL 1486 146 GL		-27.9828	8 -27.889	97 115.5613	B1 64	
1486 146 GL 1487 146 GL		-26.2948	88 -30.936	94 118.57642	B1 64	
1488 147 - LY		-21.2104 -26.3694	7 -30.918.	39 117.77294 26-119.82371	B1 64	
1489 147 LY	s H	-25.6524	9 -31.2069	9 120.47663	B1 65	
1490 147 LYS		-27.3546	3 -32.4661	4 120.21964	B1 65	
1491 147 LYS		-28.2725	832.0085	1 121.37895	B1 65	
1492 147 LYS		-29.3471	6 -33.0272	6 121.84203	B1 65	
1493 147 LYS		-28.8867	4 -34.4917	2 121.97244	B1 (5	0.00000
1495 147 LYS		-29.9661	B -35.4678	5 121.47444	B1 65	0.00000
1496 147 LYS		-30 0488	1 -30.8103 5 -37 4222	4 121.26907 3 120.71767		0.00000
1497 147 LYS		-28 53261	0 -31.4222 2 -36 7685	8 120.71767	B1 65 B1 65	0.00000
1498 147 LYS	H23	-29.15199	-37.2889	2 122.15073	B1 65	0.00000 0.00000
1499 147 LYS		-28.12445	-33.1268	9 119.09340	B1 65	0.00000
1500 147 LYS	0	-29.30235	-32.9017	4 118.83883	B1 65	0.00000
1501 148 ASP	N ·	-27.34620	-34.0091	5 118.45822	B1 66	0.00000
1502 148 ASP	H ·	-26.39747	-34.1449	118.75300	B1 66	0.00000
1503 148 ASP 1504 148 ASP	CA :	-21.19510 -27 98927	-35.0023(5 117.48362 3 118.17059	B1 66	0.00000
1505 148 ASE	CG -	-26.79528	-36.57930	119.21585	Bl 66 Bl 66	0.00000 0.00000
1506 148 ASP	OD1 -	-25.69280	-36.06325	119.05193	B1 66	0.00000
1507 148 ASP	OD2 -	-27.07650	-37.20742	120.23524	B1 66	0.00000
1508 148 ASP 1509 148 ASP	C _ :	-29.08887	-34.63043	116.79645	B1 66	0.00000
1509 148 ASP 1510 149 LEU	0 -	30.17136	-35.14808	117.04951	B1 66	0.00000
1511 149 LEU	N -	20.32399	-33.61840	115.93941	B1 67	0.00000
1512 149 LEU	•	30.07076	-32.84836	115.45008	B1 67 B1 67	0.00000 0.00000
1513 149 LEŲ		29.45399	-31.74267	114.57360	B1 67	0.00000
1514 149 LEU	CG -	30.29432	-30.58223	114.02475	31 67	0.00000
1515 149 LEU	ב נפס	30.85820	-30.92475	112.65290	31 67	0.0000
1516 149 LEU 1517 149 LEU	CD2 -	31.34761	-30.09615	115.02072	31 67	0.0000
1517 149 LEU 1518 149 LEU	C -:	31.17667	-33.69413	114.80952 1 114.83807 E	67	0.00000
1519 150 LEU	N -	30.73118	-34 86138	114.83807 E	31 67 31 68	0.00000 0.00000
1520 150 LEU	н -:	29.75579	-34.94139	114.11621 E	1 68	0.00000
1521 150 LEU	CA -	31.59782	-36.02822	114.12850 B	1 68	0.00000
1522 150 LEU	CB -3	30.74740	-37.29867	114.15286 B	1 68	0.0000
1523 150 LEU 1524 150 LEU	CG -2	9.89363	-37.44772	112.89569 B	1 68	0.00000
1524 150 LEU 1525 150 LEU	CD1 -2 CD2 -3	. 04060 ·	-38.26626	113.18796 B	1 68	0.00000
1526 150 LEU	-	7 74677 C	-38.01712	111.73915 B 115.10785 B		0.00000
1527 150 LEU		3.89001 -	-36.1/24/ -36.01901	115.10785 B	1 65 1 63	0.00000 0.00000
1528 151 GLU	к -3	2.47441 .	-36,43576	116.39428 B	1 69	0.00000
1529 151 GLU	н -3	1.52943 -	-36.57284	116.72119 B	1 65	0.00000
	CA -3	3.61295 -	-36.59512	117.30950 E	2 69	0.00000
	CB -3	3.19489 -	-36.98331	118.72928 B	1 69	0.00000
1332 151 GTA	CG -3	2.69081 -	-55.41906	118.86324 9	2 69	0.00000

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3522 344		
-31.1328/ -38.44464 118 70467 P	69	0.0000
1535 151 GLU OE2 -30.51522 -38 56462 310 20440 B	L 69	0.00000
1536 151 GLU C -34.55067 -35 41168 117 42857 PS	69	0.00000
1337 151 GLU O -35.75790 -35.57313 117.58647 BY	69	0.00000
1338 132 GLN N -33.98601 -34.19780 117.35586 BI	70	0.00000
	. 70	0.0000
1540 152 GLN CA -34.89656 -33.04752 117.40695 B1 1541 152 GLN CB -34.15680 -31.71646 117.32018 B1	70	0.00000
1542 152 GLN CG -33.46159 -31.26900 118.60094 B1	70 70	0.00000 0.00000
1543 152 GLN CD -33.23376 -29.77424 118.49387 B1	70	0.00000
1544 152 GLN OE1 -32.13068 -29.27019 118.36008 B1	70	0.00000
1545 152 GLN NE2 -34.34656 -29.04662 118.54711 B1 1546 152 GLN HE21 -35.24607 -29.46511 118.65663 B1	70	0.00000
1547 152 GLN HE22 -34.28188 -28.05358 118.46911 B1	70	0.00000
1345 132 GLN C -35.89575 -33.04746 116 27182 81	70 70	0.00000
1349 152 GLN O -37.09756 -32.85841 116.43607 Bl	70	0.00000
1330 153 ARG N -35.34435 -33.30759 115.08552 81	71	0.00000
31.5555 -53.46/09 114.49/99 RI	71	0.00000
50.23633 -33.42129 119.95927 RI	71	0.00000
1553 153 ARG CB -35.46322 -33.43707 112.66564 B1 1554 153 ARG CG -34.84280 -32.07791 112.35054 B1	71	0.00000
1555 153 ARG CD -33.884.9932 22570 131 17879 PM	71 71	0.00000
133 ARG NE -33.53171 -30.95306 110 66366 pt	71	0.00000 0.00000
133 ARG HE -33.79858 -30.09360 110.98860 B1	71	0.00000
	71	0.00000
	71	0.00000
1560 153 ARG HH11 -32.81618 -30.04802 107.55422 B1 1561 153 ARG HH12 -33.14801 -29.01024 108.90136 B1	71	0.00000
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-1, 133 W -	73	0.00000
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	73	0.00000
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595 155 ARG HH12 -42.41618 -31.75018 122.77039 31	73 73	0.00000 0.00000
	7.7	0.00000

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16	يد 162 . 61	RC IIE	50 /96	94 -40.884	55 114.1109	7 R) E	0 5.00000
2,6	62 162 AJ	RG C2			07 114.8060		0.00000
٠٠.		RG NH1			64 114.4673		0.00000
160			-50.119	79 -44.314	03 114.66150) B1 8	0.00000
160					85 113.9602		
166					63 115.3832		
166				34 -43.727			
166					88 115.64161 06 113.42487		
167					32 112.81658		
167		_			62 114.60088		
167					35 115.03933		
167	3 163 HI	S CA			57 115.23467		
167					15 116.66134		0.00000
167					9 117.55950		
167					2 118.12240		
167 167			-52.4283	0 -35.6743 4 -36.2003	31 118.03242 37 117.93420	B1 81	
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168					8 114.44466		
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1688					9 114.02377		
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2690					3 113.14586		0.00000 0.00000
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1696 1697	165 TYR 165 TYR				110.10401		0.00000
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1723	158 GLY				111.38362 E 109.94972 B		0.00000
.724	158 GLY				109.11507 B		0.00000
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172		Y O	58.026	49 -31.280	44 108.61617	7 Bl	86 0.00000
172	6 169 G <u>1</u>	א ט.	56.711	34 -33.071	69 109.00665	Bl	87 0.00000
17.		א ט.			85 109.43169		87 0.00000
1721 1729		U CA			67 108.30465		87 0.00000
1730					55 108.28846 91 107.43481		87 0.00000 87 0.00000
1732					19 107.43401		87 0.00000
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1739					1 112.35798		88 0.00000
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1755	171 PHE	٥.			7 108.07691		0.00000
1756 1757	172 THR 172 THR	И, Н			107.69903 108.09636		0.00000
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	174 GLN 174 GLN	CA CB			102.94243 8		
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	175 ARG				100.50018 B		
1788	175 ARG	CB .	-61.85817	-30.36631	100.16958 B	1 93	0.00000

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987ג			-61.2439	5 -30.2691	4 98.77310 B1	93	0.00000
1790		G CD	-60.9159	7 -28.0420	9 98.34709 B1	93	0.00000
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	178 HIS		66.60728		96.05191 Bl	96 ·	0.00000
	178 HIS		62.34621		97.09131 B1	96	0.00000
	178 HIS		61.22615		97.07091 B1	96	0.00000
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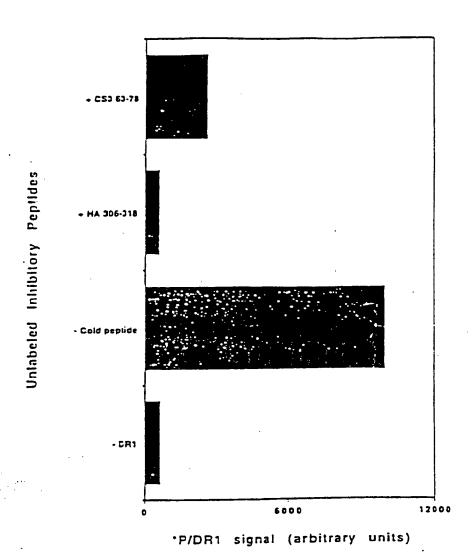
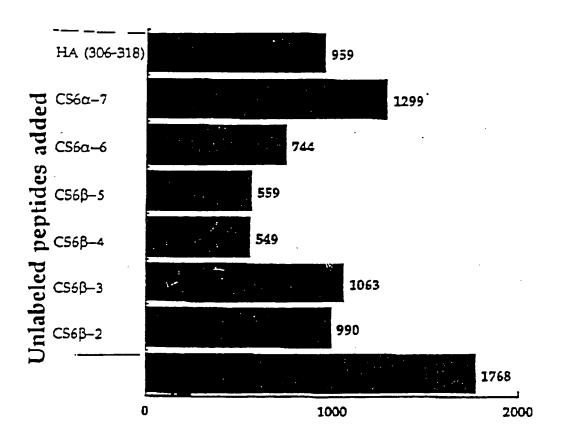


FIG. 31

Inhibition of 125 I HA (306-318)/DRI by unlabeled CSG of and B peptides



*HA/DR1 compact dimer signal (densitometric units)

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

A. CLASSIFICATION OF SUBJECT MATTER								
IPC(5) :A61K 39/00, 39/02, 39/12, 37/02, 35/14 US CL :424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75								
According to International Patent Classification (IPC) or to both national classification and IPC								
B. FIELDS SEARCHED								
Minimum documentation searched (classification system	followed by classification symbols)							
U.S. : 424/185.1, 186.1, 190.1, 242.1; 530/327, 326, 333, 334, 388.75								
Documentation searched other than minimum documentation	tion to the extent that such documents are included in the fields searched							
Electronic data base consulted during the international s	cearch (name of data base and, where practicable, search terms used)							
C. DOCUMENTS CONSIDERED TO BE RELEVANT								
Category* Citation of document, with indication,	where appropriate, of the relevant passages Relevant to claim No.							
X The Journal of Immunology	, Volume 150, No. 8, Part II, 1, 3-20							
issued 15 April 1993, Naus	s et al., " Binding Interactions of							
Peptides in a Structural Hon	nology Model of the DR1 Class							
MHC ", page 41A, Abstract	221, see entire abstract.							
X Nature, Volume 358, issued	27 August 1992, Chicz et al., 12							
	cessed Peptides Bound to HLA-							
	HC-related Molecule and are 1, 3-7							
	s 764-768, see page 766, Table							
2, and Table 3.								
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X Further documents are listed in the continuation of	of Box C. See patent family annex.							
Special categories of cited documents:	"T" later document published after the international filing date or priority							
"A" document defining the general state of the art which is not co	maidered date and not in conflict with the application but cited to understand the principle or theory underlying the invention							
"E" earlier document published on or after the international filing	document of particular relevance; the claimed invention cannot be considered novel or cannot be considered to involve an inventive step							
"L" document which may throw doubts on priority claim(a) or cited to establish the publication date of another citation	which is when the document is taken alone or other							
special reason (as specified) "O" document referring to an oral disclosure, use, exhibition means	"Y" document of particular relevance; the claimed invention cannot be considered to involve an inventive step when the document is or other combined with one or more other such documents, such combination being obvious to a person skilled in the art							
P document published prior to the international filing date but is the priority date claimed	• • • • • •							
Date of the actual completion of the international search Date of mailing of the international search report								
01 SEPTEMBER 1994	1 3 SEP 1994.							
Name and mailing address of the ISA/US Authorized officer								
Commissioner of Patents and Trademarks Box PCT Washington D. C. 2002	H. Sidberry W. Kuza fa							
Washington, D.C. 20231 Facsimile No. (703) 305-3230	Telephone No. (703) 308-0196							
Form PCT/ISA/210 (second sheet)(July 1992)*								

INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

		C 17US9470369	• •
C (Continu	uion). DOCUMENTS CONSIDERED TO BE RELEVANT		
Category*	Citation of document, with indication, where appropriate, of the relevant	Relevant to claim No	
Y	The Journal of Immunology, Volume 150, No. 2, issued January 1993, Boehncke et al., "The Importance of Dom Negative Effects of Amino Acid Side Chain Substitution Peptide-MHC Molecule Interactions and T Cell Recogniti pages 331-341, see Abstract, on page 331.	8-11	
x	The EMBO Journal, Volume 9, No. 6, issued 1990, Jardal., "Peptide binding to HLA-DR1: a Peptide with most r substituted to alanine retains MHC binding", pages 1797-page 1798, page 1800, figure 4, and page 1801, figure 7.	512	
Ý	Nature, Volume 332, issued 28 April 1988, Brown et al., hypothetical model of the foreign antigen binding site of (histocompatibility molecules", pages 845-850, see pages 8	Class II	1, 3, 4
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INTERNATIONAL SEARCH REPORT

International application No. PCT/US94/05697

Box 1 Observations where certain claims were found unsearchable (Continuation of item 1 of first sheet)									
This international report has not been established in respect of certain claims under Article 17(2)(a) for the following reasons:									
Claims Nos.: 2 because they relate to subject matter not required to be searched by this Authority, namely:									
Claim 2 is directed to a computerized model which encompasses scientific theory and computer programs to the extent that the International Searching Authority is not equipped to search prior art concerning such programs. Accordingly claim 2 is withdrawn from search under PCT Rule 39 and PCT Article 17(2)(a)(i).									
Claims Nos.: because they relate to parts of the international application that do not comply with the prescribed requirements to such an extent that no meaningful international search can be carried out, specifically:									
Claims Nos.: because they are dependent claims and are not drafted in accordance with the second and third sentences of Rule 6.4(a).									
Box II Observations where unity of invention is lacking (Continuation of item 2 of first sheet)									
This International Searching Authority found multiple inventions in this international application, as follows:									
1. As all required additional search fees were timely paid by the applicant, this international search report covers all searchable claims.									
2. As all searchable claims could be searched without effort justifying an additional fee, this Authority did not invite payment of any additional fee.									
3. As only some of the required additional search fees were timely paid by the applicant, this international search report covers only those claims for which fees were paid, specifically claims Nos.:									
4. No required additional search fees were timely paid by the applicant. Consequently, this international search report is restricted to the invention first mentioned in the claims; it is covered by claims Nos.:									
Remark on Protest									
No protest accompanied the payment of additional search fees.									